10/018463 531 Rec'd PCT/PTC 19 DEC 2001

IN THE UNITED STATES PATENT AND TRADEMARK OFFICE REQUEST FOR FILING NATIONAL PHASE OF

PCT APPLICATION UNDER 35 U.S.C. 371 AND 37 CFR 1.494 OR 1.495

To:

Hon. Commissioner of Patents

Washington, D.C. 20231

**

TRANS	MITTAL LETTER TO THE UNITED S	ES	Atty Dkt:			46 /160					
	NATED/ELECTED OFFICE (DO/EO/U			,				<u>M#</u>	/Client Ref	•	
From:	Manelli Dension & Selter:			Date:	:	Decembe	er 19,	2001			
	This is a REQUEST for FILING a PO	CT/US	SA Natio	nal Phase A	4pr	olication ba	sed (on:			
1. Li	International Application	2.	Interna	ational Filing	j D	ate	3.	Earlies	st Priority Date	e Claimed	
r- C3	PCT/JP00/04136		23	June	2	2000	_	24	June	1999	
	û country code		Day	MONTH		Year		Day use itei	MONTH m 2 if no earli	Year er priority)	
	Measured from the earliest priority date in item 3, this PCT/USA National Phase Application Request is being filed within:										
	(a) 20 months from above item 3	date	(b)	☐ 30 month	hs	from abov	e iten	n 3 date	е,		
	(c) Therefore, the due date (unexter	<u>idable</u>	e) is _2	4 December	r 2	001					
	Title of Invention: <u>HETEROCYCLIC DICARBOXYLIC ACID DIAMIDE DERIVATIVES</u> , <u>AGRICULTURAL AND HORTICULTURAL INSECTICIDES AND METHOD OF USING THE SAME</u>										
a bar .	Inventor(s) Takeshi Katsuhira, Takashi Furuya, Makoto Gotoh, Masanori Tohnishi, Hideo Takaishi, Kazuyuki Sakata, Masayuki Morimoto and Akira Seo pplicant herewith submits the following under 35 U.S.C. 371 to effect filing:										
7.							371 (f)).			
8.	☐ A copy of the International Application as filed (35 U.S.C. 371(c)(2)) is transmitted herewith (file if in English but, if in foreign language, file only if not transmitted to PTO by the International Bureau) including:										
	 a. Request; b. Abstract; c. pgs. Spec. and Claims; d. sheet(s) Drawing which are informal formal of size A4 11" 										
9.	□ A copy of the International Ap						ne Int	ernatio	onal Bureau.		
				•							
10.	A translation of the International a. is transmitted herewith in (3) 189 pgs. Spec. (4) 0 sheet(s) Dr	ncludi and C awing	ng: (1) Claims; g which	∐ Request	t; (2) <u>⊠</u> Abst	ract;				
	 b. is not required, as the application was filed in English. c. is not herewith, but will be filed when required by the forthcoming PTO Missing Requirements Notice per Rule 494(c) if box 4(a) is X'd or Rule 495(c) if box 4(b) is X'd. d. Translation verification attached (not required now). 										

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Page 2 of 3

RE: USA National Fifing of PCT/JP00/04136

	24.	Atta	ched:	1. Form PCT/IB 2. Cover page of		d international a	ກວີໄດ້ເ application WC		ntains Er	nglish Abs	tract)
	25.	Prel	imina	ry Amendment:	ATTACHE		iminary Amen Preliminary An		ior to fee	e calculati	on)
	25.5	Per	ltem 1	7.c2, <u>cancel origi</u>	nal pages	#, claims	#, Drav	ving Sheets #_			
	26. Based	6. Calculation of the U.S. National Fee (35 U.S.C. 371 (c)(1)) and other fees is as follows: ased on amended claim(s) per above item(s) 12, 14, 17, 25, 25, 25.5 (hilite)									
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	27.							SUBTOTAL =	\$890	0.00	=
	28.	If As	signm	ent box 19 above	is X'd, add	X'd, add Assignment Recording fee of\$40					(581)
	29. Attached is a check to cover the TOTAL FEES \$88							\$890	0.00	=	
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	filed, or whi or hereafter duplicate co	ch shou relative opy of th	ild have be to this ap his sheet is	e Commissioner is hereby a een filed herewith or concer oplication and the resulting (s attached. I does not authorize charg	ning any paper fi Official document	rge any fee specifically led hereafter, and whic under Rule 20, or crec	h may be required un lit any overpayment, t	der Rules 16-18 and 49 to our Account/Order N	92 (missing o	r insufficient fe	e only) now
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By Atty:					By Atty:	Paul E. White	Re	Reg. No. 32,011			
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IN THE UNITED STATES PATENT AND TRADEMARK OFFICE

#4/ B

In re PATENT APPLICATION of KATSUHIRA, et al.

Group Art Unit: Not Assigned

Appln. No.: Not Assigned

Examiner: Not Assigned

Filed: December 3, 2001

International Appln. No. PCT/JP00/04136

Title: HETEROCYCLIC DICARBOXYLIC ACID DIAMIDE DERIVATIVES,

AGRICULTURAL AND HORTICULTURAL INSECTICIDES AND

METHOD OF USING THE SAME

December 19, 2001

SECOND PRELIMINARY AMENDMENT: TO BE ENTERED AFTER CALCULATION OF FILING FEE

Hon. Commissioner of Patents and Trademarks Washington, D.C. 20231

Sir:

Please enter the following Preliminary Amendment of the subject new application prior to calculation of the fee for filing the application.

IN THE SPECIFICATION (see the attached Appendix for the changes made to effect the below amendments to the specification):

Paragraph at page 157, line 27 to page 158, line 1

Physical property:

compound No. 1382 m.p. 164-166 . Yield 36%.

compound No. 1414 m.p. 167-169 . Yield 36%.

Paragraph at page 163 starting at line 12

As a result, the following compounds were found to have a corrected mortality of 90% or more: compound Nos. 13, 18, 32, 54, 55, 57, 127, 136, 230, 242, 258, 464, 484, 512, 524, 737, 785, 794, 795, 804, 805, 821, 989, 990, 1009, 1095, 1110, 1127, 1158, 1189, 1204, 1220, 1221, 1247, 1249, 1251, 1255, 1267, 1269, 1271, 1275, 1303, 1306, 1313, 1320, 1414, 1429, 1473, 1505, 1548 and 1549.

IN THE CLAIMS:

Please amend claim 1 as follows (see the attached Appendix for the changes made to effect the below claim):

1. A heterocyclic dicarboxylic acid diamide derivative represented by the general formula (I):

$$\begin{array}{c} Xn \\ \text{Het} \\ & \\ Z^2 \end{array} \xrightarrow{NR^1R^2} Ym \\ & \\ Z^2 \xrightarrow{R_3} \xrightarrow{B^4-B^3} Ym \end{array}$$

{wherein R^1 , R^2 and R^3 , which may be the same or different, are hydrogen atoms, (C_3-C_6) cycloalkyl groups, halo (C_3-C_6) cycloalkyl groups or $-A^1-(R^4)$ r (wherein A^1 is a (C_1-C_8) alkylene group, a (C_3-C_6) alkenylene group or a (C_3-C_6) alkynylene group, R^4 , which may be the same or different, are hydrogen atoms; halogen atoms; cyano groups; nitro groups; halo (C_1-C_6) alkyl groups; (C_3-C_6) cycloalkyl groups; halo (C_3-C_6) cycloalkyl groups;

C₆)cycloalkyl groups; (C₁-C₆)alkoxycarbonyl groups; di(C₁-C₆)alkoxyphosphoryl groups whose (C₁-C₆)alkoxy groups may be the same or different; di(C₁-C₆)alkoxythiophosphoryl groups whose (C₁-C₆)alkoxy groups may be the same or different; diphenylphosphino groups; diphenylphosphono groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^2-R^5$ (wherein A^2 is $-O_-$, $-S_-$, -SO₂--, -N(R⁶)- (wherein R⁶ is a hydrogen atom; a (C₁-C₆)alkylcarbonyl group; a halo(C_1 - C_6)alkylcarbonyl group; a (C_1 - C_6)alkoxycarbonyl group; a phenylcarbonyl group; a substituted phenylcarbonyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups. halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenyl(C₁-C₄)alkoxycarbonyl group; a substituted phenyl(C₁-C₄)alkoxycarbonyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C1-

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 C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁- C_6)alkylsulfonyl groups; a (C_1 - C_6)alkylsulfonyl group; or a halo(C_1 - C_6)alkylsulfonyl group), -C(=0) or $-C(=NOR^7)$ (wherein R^7 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo(C_1 - C_6)alkyl group; a (C_3 - C_6)alkenyl group; a halo(C_3 - C_6)alkenyl group; a (C₃-C₆)alkynyl group; a cyclo(C₃-C₆)alkyl group; a phenyl(C₁-C₄)alkyl group; or a substituted phenyl(C₁-C₄)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C₁- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups), and R⁵ is a hydrogen atom; a (C₁-C₆)alkyl group; a halo(C₁- C_6)alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) alkenyl group; a (C_3-C_6) alkynyl group; a halo(C_3 - C_6)alkynyl group; a (C_3 - C_6)cycloalkyl group; a halo(C_3 - C_6)cycloalkyl group; a (C_1-C_6) alkoxy (C_1-C_6) alkyl group; a (C_1-C_6) alkylthio (C_1-C_6) alkyl group; a formyl group; a (C₁-C₆)alkylcarbonyl group; a halo(C₁-C₆)alkylcarbonyl group; a (C₁-C₆)alkoxycarbonyl group; a mono(C₁-C₆)alkylaminocarbonyl group; a di(C₁-C₆)alkylaminocarbonyl group whose (C₁-C₆)alkyl groups may be the same or different; a mono(C₁-C₆)alkylaminothiocarbonyl group; a di(C₁-C₆)alkylaminothiocarbonyl group whose (C₁-C₆)alkyl groups may be the same or different; a di(C₁-C₆)alkoxyphosphoryl group whose (C₁-C₆)alkoxy groups may be the same or different; a di(C₁-C₆)alkoxythiophosphoryl group whose (C₁-C₆)alkoxy groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are

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selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁- C_6)alkylthio groups, (C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C_1 -C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenyl(C₁-C₄)alkyl group; a substituted phenyl(C₁-C₄)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁- C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups), and r is an integer of 1 to 4),

provided that R¹ and R² are not hydrogen atoms at the same time,

R¹ and R² may form a 4 to 7 membered ring by combining to each
other, in which the ring may contain the same or different 1 to 3 hetero atoms
selected from the group consisting of oxygen atom, sulfur atom and nitrogen atom.

Het is a heterocyclic group represented by any of the following formulas Q1 to Q22:

$$Q1 = \begin{cases} x_{1} & x_{1}$$

(wherein X, which may be the same or different, are halogen atoms; cyano groups; nitro groups; (C₃-C₆)cycloalkyl groups; halo(C₃-C₆)cycloalkyl groups; tri(C₁-C₆)alkylsilyl groups whose (C₁-C₆)alkyl groups may be the same or different: phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1 - C_6)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or -A³-R⁸ [wherein A³ is -O-, -S-, -SO-, - SO_{2} --, $-N(R^6)$ - (wherein R^6 is as defined above), -C(=O)-, $-C(=NOR^7)$ -(wherein R⁷ is as defined above), a (C₁-C₆)alkylene group, a halo(C₁-C₆)alkylene group, a (C₂-C₆)alkenylene group, a halo(C₂-C₆)alkenylene group, a (C₂- C_6)alkynylene group or a halo(C_3 - C_6)alkynylene group, and R^8 is as follows: (1) when A^3 is $-O_-$, $-S_-$, $-S_-$, $-S_-$ or $-N(R^6)_-$ (wherein R^6 is as defined above). then R⁸ is a halo(C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkenyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl

groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or $-A^4-R^9$ (wherein A^4 is a (C₁-C₆)alkylene group, a halo(C₁-C₆)alkylene group, a (C₃-C₆)alkenylene group, a halo(C₃-C₆)alkenylene group, a (C₃-C₆)alkynylene group or a halo(C₃-C₆)alkynylene group, and R⁹ is a hydrogen atom; a halogen atom; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or -A⁵-R¹⁰ (wherein A⁵ is $-O_{-}$, $-S_{-}$, $-SO_{-}$, $-SO_{2}$ or -C(=O), and R^{10} is a $(C_{1}-C_{6})$ alkyl group; a halo $(C_{1}-C_{6})$ C₆)alkyl group; a (C₃-C₆)alkenyl group; a halo(C₃-C₆)alkenyl group; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups)),

(2) when A^3 is -C(=0)— or $-C(=NOR^7)$ — (wherein R^7 is as defined above), then R^8 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo (C_1-C_6) alkyl group; a (C_2-C_6) alkenyl group; a halo(C₂-C₆)alkenyl group; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C_1-C_6) alkoxy group; a (C_1-C_6) alkylthio group; a mono (C_1-C_6) alkylamino group; a di(C₁-C₆)alkylamino group whose (C₁-C₆)alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenylamino group; a substituted phenylamino group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁- C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl

groups, and

(3) when A^3 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6) alkylene group, a (C_3-C_6) alkylene C_6)alkenylene group, a halo(C_2 - C_6)alkenylene group, a (C_2 - C_6)alkynylene group or a halo(C₃-C₆)alkynylene group, then R⁸ is a hydrogen atom; a halogen atom; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a tri(C₁-C₆)alkylsilyl group whose (C₁-C₆)alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or -A⁶-R¹¹ (wherein A⁶ is -O-, -S-, -SO- or -SO₂--, and R¹¹ is a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen

atoms. (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁- C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthi C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or $-A^7-R^{12}$ (wherein A^7 is a (C₁-C₆)alkylene group, a halo(C₁-C₆)alkylene group, a (C₂-C₆)alkenylene group, a halo(C₂- C_6)alkenylene group, a (C_2 - C_6)alkynylene group or a halo(C_3 - C_6)alkynylene group, and R¹² is a hydrogen atom; a halogen atom; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C₁-C₆)alkoxy group; a halo(C₁-C₆)alkoxy group; a (C₁-C₆)alkylthio group; a halo(C₁-C₆)alkylthio group; a (C₁-C₆)alkylsulfinyl group; a halo(C₁-C₆)alkylsulfinyl group; a (C₁-C₆)alkylsulfonyl group; a halo(C₁-C₆)alkylsulfonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C1- C_6)alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenoxy group; a substituted phenoxy group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenylthio group; a substituted phenylthio group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-

 C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkylthio groups, halo(C_1 - C_6)alkylthio groups, halo(C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups)], and n is an integer of 0 to 3,

X may form a condensed ring by combining together with the adjacent atoms in the heterocyclic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C₁-C₆)alkyl groups; halo(C₁-C₆)alkyl groups; (C₁-C₆)alkoxy groups; halo(C₁-C₆)alkoxy groups; (C₁-C₆)alkylthio groups; halo(C₁-C₆)alkylthio groups; (C₁- C_6)alkylsulfinyl groups; halo(C_1 - C_6)alkylsulfinyl groups; (C_1 - C_6)alkylsulfonyl groups; halo(C₁-C₆)alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1 - C_6)alkylthio groups, (C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups,

W is O, S or N-R¹³ (wherein R¹³ is a (C₁-C₆)alkyl group; a halo(C₁-

 C_6)alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) alkenyl group; a (C_3-C_6) alkynyl group; a halo (C_3-C_6) alkynyl group; a (C_1-C_6) alkoxy group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, a phenyl (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups are phenyl (C_1-C_6) alkyl group; or a substituted phenyl (C_1-C_6) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfonyl groups,

B¹, B², B³ and B⁴, which may be the same or different, are carbon atoms or nitrogen atoms,

Y, which may be the same or different, are halogen atoms; cyano groups; nitro groups; halo(C_3 - C_6)cycloalkyl groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkylthio groups, halo(C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo($C_$

 C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ (wherein A^3 and R^8 are as defined above), and m is an integer of 1 to 5,

Y may form a condensed ring by combining together with the adjacent carbon atoms in the aromatic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; (C_1-C_6) alkoxy groups; halo (C_1-C_6) alkoxy groups; halo (C_1-C_6) alkyl groups; halo (C_1-C_6) C₆)alkoxy groups; (C₁-C₆)alkylthio groups; halo(C₁-C₆)alkylthio groups; (C₁-C₆)alkylsulfinyl groups; halo(C₁-C₆)alkylsulfinyl groups; (C₁-C₆)alkylsulfonyl groups; halo(C₁-C₆)alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups, and each of Z^1 and Z^2 is an oxygen atom or a sulfur atom.

provided that:

(1) when Het is Q2, Q6, Q7 or Q9 and B1, B2, B3 and B4 are carbon atoms at the

same time, then Ym is other than 3-chloro-2-methyl group, 3-chloro-2,6-diethyl group, 5-chloro-2-methyl group, 2,6-diethyl group, 4-chloro-2-fluoro group and 2-ethyl-6-methyl group,

- (2) when Het is Q4 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 2,5-dichloro group, 2,4-difluoro group, 2,6-difluoro group, 3-chloro-2-methyl group, 5-chloro-2-methyl group, 5-fluoro-2-methyl group, 2,6-dimethyl group, 2,6-diethyl group, 2-ethyl-6-methyl group, 2-methoxy-5-nitro group, 2-methoxy-5-methyl group, 2,6-diethoxy group, 3-bromo-2-methyl group, 3-fluoro-2-methyl group, 3-iodo-2-methyl group, 3-cyano-2-methyl group, 3-difluoromethoxy-2-methyl group, 5-chloro-2-ethyl group, 2,5-dimethyl group, 2,3-dichloro group, 3-chloro-2,6-diethyl group, 4-trifluoromethyl group, 3-methoxycarbonyl-2-methyl group, 3-trifluoromethyl-2-methyl group, 3,5-dichloro-2,6-diethyl group, 3,4-dichloro group, 3-(methoxycarbonylmethyloxy)-2-methyl group, 2-methyl-3-nitro group and 4-trifluoromethoxy group,
- (3) when Het is Q9, R² and R³ are hydrogen atoms at the same time, Xn is a 2-phenyl group, R¹ is a n-propyl group or an i-propyl group and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 4-pentafluoroethyl-2-methyl group, and
- (4) when Het is Q10 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 5-chloro-2-methyl group, 5-fluoro-2-methyl group, 2,5-dimethyl group and 2,6-diethyl group, and
- (5) when Het is Q10 and B¹, B², B³ and B⁴ are carbon atoms at the same time, Xn is other than 5,6-dimethyl group }.

REMARKS

This Preliminary Amendment revises claim 1 in accordance with the Amendment under PCT Article 34 that was made in the Japanese language during the international phase of the subject international application.

The specification has been amended as shown above to correct typographical errors that were made in producing the presently filed English translation from the original application filed in Japanese.

No new matter has been added.

Entry of this amendment and favorable consideration of this application are respectfully requested.

Respectfully submitted,

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APPENDIX SHOWING REVISIONS OF SPECIFICATION AND CLAIMS

Proposed Amendments To The Specification Showing Deletions And Insertions

Paragraph at page 157, line 27 to page 158, line 1

Physical property:

compound No. [1381] <u>1382</u> m.p. 164–166 . Yield 36%. compound No. [1412] <u>1414</u> m.p. 167–169 . Yield 36%.

Paragraph at page 163 starting at line 12.

As a result, the following compounds were found to have a corrected mortality of 90% or more: compound Nos. 13, 18, 32, 54, 55, 57, 127, 136, 230, 242, 258, 464, 484, 512, 524, 737, 785, 794, 795, 804, 805, 821, 989, 990, 1009, 1095, 1110, 1127, 1158, 1189, 1204, 1220, 1221, 1247, 1249, 1251, 1255, 1267, 1269, 1271, 1275, 1303, 1306, 1313, 1320, 1414, 1429, 1473, 1505, 1548 and [1540] 1549.

Proposed Amendments To Claim 1 Showing Deletions And Insertions.

1. A heterocyclic dicarboxylic acid diamide derivative represented by the general formula (I):

$$\begin{array}{c|c} Xn & Z^1 \\ & NR^1R^2 \\ & & B^1 = B^2 \\ & & N \end{array}$$

$$Xn & NR^1R^2 \\ & & NR^1R^2$$

{wherein R¹, R² and R³, which may be the same or different, are hydrogen atoms, (C₃-C₆)cycloalkyl groups, halo(C₃-C₆)cycloalkyl groups or -A¹-(R⁴)r (wherein A¹ is a (C_1-C_8) alkylene group, a (C_3-C_6) alkenylene group or a (C_3-C_6) alkynylene group, \mathbb{R}^4 . which may be the same or different, are hydrogen atoms; halogen atoms; cyano groups; nitro groups; halo(C₁-C₆)alkyl groups; (C₃-C₆)cycloalkyl groups; halo(C₃-C₆)cycloalkyl groups; (C₁-C₆)alkoxycarbonyl groups; di(C₁-C₆)alkoxyphosphoryl groups whose (C₁-C₆)alkoxy groups may be the same or different; di(C₁- C_6)alkoxythiophosphoryl groups whose (C_1 - C_6)alkoxy groups may be the same or different; diphenylphosphino groups; diphenylphosphono groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen. atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkyl groups, halo (C_1-C_6) C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^2-R^5$ (wherein A^2 is $-O_-$, $-S_-$, $-SO_-$, -SO₂--, -N(R⁶)- (wherein R⁶ is a hydrogen atom; a (C₁-C₆)alkylcarbonyl group; a

halo(C₁-C₆)alkylcarbonyl group; a (C₁-C₆)alkoxycarbonyl group; a phenylcarbonyl group; a substituted phenylcarbonyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups. halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁- C_6)alkylsulfonyl groups; a phenyl(C_1 - C_4)alkoxycarbonyl group; a substituted phenyl(C_1 - C_4)alkoxycarbonyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms. (C₁- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁- C_6)alkylsulfonyl groups; a (C_1-C_6) alkylsulfonyl group; or a halo (C_1-C_6) alkylsulfonyl group), -C(=O)— or $-C(=NOR^7)$ — (wherein R^7 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo(C₁-C₆)alkyl group; a (C₃-C₆)alkenyl group; a halo(C₃-C₆)alkenyl group; a (C_3-C_6) alkynyl group; a cyclo (C_3-C_6) alkyl group; a phenyl (C_1-C_4) alkyl group; or a substituted phenyl(C₁-C₄)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms. (C₁- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups), and R⁵ is a hydrogen atom; a (C₁-C₆)alkyl group; a halo(C₁-C₆)alkyl group; a (C₃-C₆)alkenyl group; a halo(C₃-C₆)alkenyl group; a (C₃-C₆)alkynyl group; a halo(C₃-C₆)alkynyl group; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C_1-C_6) alkoxy (C_1-C_6) alkyl group; a (C_1-C_6) alkyl group; a

formyl group; a (C₁-C₆)alkylcarbonyl group; a halo(C₁-C₆)alkylcarbonyl group; a (C₁-C₆)alkoxycarbonyl group; a mono(C₁-C₆)alkylaminocarbonyl group; a di(C₁-C₆)alkylaminocarbonyl group whose (C₁-C₆)alkyl groups may be the same or different; a mono(C₁-C₆)alkylaminothiocarbonyl group; a di(C₁-C₆)alkylaminothiocarbonyl group whose (C₁-C₆)alkyl groups may be the same or different; a di(C₁-C₆)alkoxyphosphoryl group whose (C₁-C₆)alkoxy groups may be the same or different; a di(C₁-C₆)alkoxythiophosphoryl group whose (C₁-C₆)alkoxy groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenyl(C₁-C₄)alkyl group; a substituted phenyl(C₁-C₄)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups), and r is an integer of 1 to 4),

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provided that R^1 and R^2 are not hydrogen atoms at the same time, R^1 and R^2 may form a 4 to 7 membered ring by combining to each other, in which the ring may contain the same or different 1 to 3 hetero atoms selected from the group consisting of oxygen atom, sulfur atom and nitrogen atom,

Het is a heterocyclic group represented by any of the following formulas Q1 to Q22:

$$Q1 = \begin{cases} Xn & 4 & 3 \\ 5 & 1 \\ 1 & (0) p \end{cases} \qquad Q2 = \begin{cases} Xn & 5 \\ 6 & 1 \\ 1 & 2 \end{cases} \qquad Q3 = \begin{cases} Xn & 2 \\ 1 & 1 \\ 1 & 3 \end{cases} \qquad Q4 = \begin{cases} Xn & 1 \\ 1 & 1 \\ 1 & 3 \end{cases}$$

$$Q9 = \begin{array}{c} Xn & (0) p \\ Xn & (0) q \\ & & & \\$$

$$Q21 = \begin{array}{c} Xn & 3 \\ 2N & 1 \end{array}$$

$$Q22 = \begin{array}{c} Xn & 3 \\ 2N & 1 \end{array}$$

(wherein X, which may be the same or different, are halogen atoms; cyano groups; nitro groups; (C₃-C₆)cycloalkyl groups; halo(C₃-C₆)cycloalkyl groups; tri(C₁-C₆)alkylsilyl groups whose (C₁-C₆)alkyl groups may be the same or different; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C1-C6)alkyl groups, halo(C1- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or -A³-R⁸ [wherein A³ is -O-, -S-, -SO-, - SO_2 --, $-N(R^6)$ - (wherein R^6 is as defined above), -C(=O)--. $-C(=NOR^7)$ -(wherein R⁷ is as defined above), a (C₁-C₆)alkylene group, a halo(C₁-C₆)alkylene group, a (C₂-C₆)alkenylene group, a halo(C₂-C₆)alkenylene group, a (C₂-C₆)alkynylene group or a halo(C₃-C₆)alkynylene group, and R⁸ is as follows: (1) when A^3 is $-O_{-}$, $-S_{-}$, $-S_{-}$, $-S_{-}$ or $-N(R^6)_{-}$ (wherein R^6 is as defined above). then R^8 is a halo(C_3 - C_6)cycloalkyl group; a halo(C_3 - C_6)cycloalkenyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C1-C6)alkyl groups, halo(C1- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl

groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^4-R^9$ (wherein A^4 is a (C_1 - C_6)alkylene group, a halo(C₁-C₆)alkylene group, a (C₃-C₆)alkenylene group, a halo(C₃-C₆)alkenylene group, a (C₃-C₆)alkynylene group or a halo(C₃-C₆)alkynylene group. and R9 is a hydrogen atom; a halogen atom; a (C3-C6)cycloalkyl group; a halo(C3-C₆)cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; or -A⁵-R¹⁰ (wherein A⁵ is -O-, -S-, -SO-, $-SO_2-$ or -C(=O), and R^{10} is a (C_1-C_6) alkyl group; a halo (C_1-C_6) C₆)alkyl group; a (C₃-C₆)alkenyl group; a halo(C₃-C₆)alkenyl group; a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups; (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C1-C6)alkyl groups, halo(C1 C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups)),

(2) when A^3 is -C(=O)— or $-C(=NOR^7)$ — (wherein R^7 is as defined above), then R^8 is a hydrogen atom; a (C₁-C₆)alkyl group; a halo(C₁-C₆)alkyl group; a (C₂-C₆)alkenyl group; a halo(C₂-C₆)alkenyl group; a (C₃-C₆)cycloalkyl group; a (C₁-C₆)alkoxy group; a (C₁-C₆)alkylthio group; a mono(C₁-C₆)alkylamino group; a di(C₁-C₆)alkylamino group whose (C₁-C₆)alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C1- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenylamino group; a substituted phenylamino group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁- C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1 - C_6)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl

groups, and

(3) when A^3 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6) alkylene group, a (C_3-C_6) alkylene C₆)alkenylene group, a halo(C₂-C₆)alkenylene group, a (C₂-C₆)alkynylene group or a halo(C_3 - C_6)alkynylene group, then R^8 is a hydrogen atom; a halogen atom; a (C_3 -C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a tri(C₁-C₆)alkylsilyl group whose (C₁-C₆)alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^6-R^{11}$ (wherein A^6 is -O-, -S-, -SO- or -SO₂-, and R¹¹ is a (C₃-C₆)cycloalkyl group; a halo(C₃-C₆)cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C1-C6)alkyl groups, halo(C1- C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁- C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen

atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^7-R^{12}$ (wherein A^7 is a (C_1 - C_6)alkylene group, a halo(C₁-C₆)alkylene group, a (C₂-C₆)alkenylene group, a halo(C₂-C₆)alkenylene group, a (C₂-C₆)alkynylene group or a halo(C₃-C₆)alkynylene group, and R^{12} is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a halogen atom; C₆)cycloalkyl group; a (C₁-C₆)alkoxy group; a halo(C₁-C₆)alkoxy group; a (C₁-C₆)alkylthio group; a halo(C₁-C₆)alkylthio group; a (C₁-C₆)alkylsulfinyl group; a halo(C₁-C₆)alkylsulfinyl group; a (C₁-C₆)alkylsulfonyl group; a halo(C₁-C₆)alkylsulfonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms. (C1- C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenoxy group; a substituted phenoxy group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; a phenylthio group; a substituted phenylthio group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-

 C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, (C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkylthio groups, halo(C_1 - C_6)alkylthio groups, halo(C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C_1 - C_6)alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups))], and n is an integer of 0 to 3,

X may form a condensed ring by combining together with the adjacent atoms in the heterocyclic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; (C_1-C_6) alkoxy groups; halo (C_1-C_6) alkyl groups; halog C₆)alkoxy groups; (C₁-C₆)alkylthio groups; halo(C₁-C₆)alkylthio groups; (C₁-C₆)alkylsulfinyl groups; halo(C₁-C₆)alkylsulfinyl groups; (C₁-C₆)alkylsulfonyl groups: halo(C₁-C₆)alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups. halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups. halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups,

W is O, S or N-R¹³ (wherein R¹³ is a (C₁-C₆)alkyl group; a halo(C₁-

 C_6)alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) alkenyl group; a (C_3-C_6) alkynyl group; a halo (C_3-C_6) alkynyl group; a (C_1-C_6) alkoxy group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfonyl groups; a phenyl (C_1-C_6) alkyl group; or a substituted phenyl (C_1-C_6) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfonyl groups, halo (C_1-C_6) alkyls

 $\mathsf{B}^1,\,\mathsf{B}^2,\,\mathsf{B}^3$ and $\mathsf{B}^4,$ which may be the same or different, are carbon atoms or nitrogen atoms,

Y, which may be the same or different, are halogen atoms; cyano groups; nitro groups; halo(C_3 - C_6)cycloalkyl groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkyl groups, halo(C_1 - C_6)alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1 - C_6)alkylthio groups, (C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups and halo(C_1 - C_6)alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6)alkyl groups, halo(C_1 -

 C_6)alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ (wherein A^3 and R^8 are as defined above), and m is an integer of 1 to 5,

Y may form a condensed ring by combining together with the adjacent carbon atoms in the aromatic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C₁-C₆)alkyl groups; halo(C₁-C₆)alkyl groups; (C₁-C₆)alkoxy groups; halo(C₁- C_6)alkoxy groups; (C_1 - C_6)alkylthio groups; halo(C_1 - C_6)alkylthio groups; (C_1 - C_6)alkylsulfinyl groups; halo(C_1 - C_6)alkylsulfinyl groups; (C_1 - C_6)alkylsulfonyl groups; halo(C₁-C₆)alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups. halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C₁-C₆)alkylthio groups, (C₁-C₆)alkylsulfinyl groups, halo(C₁-C₆)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆)alkyl groups, halo(C₁-C₆)alkyl groups, (C₁-C₆)alkoxy groups, halo(C₁-C₆)alkoxy groups, (C₁-C₆)alkylthio groups, halo(C_1 - C_6)alkylthio groups, (C_1 - C_6)alkylsulfinyl groups, halo(C_1 - C_6)alkylsulfinyl groups, (C₁-C₆)alkylsulfonyl groups and halo(C₁-C₆)alkylsulfonyl groups, and each of Z^1 and Z^2 is an oxygen atom or a sulfur atom.

provided that:

(1) when Het is Q2, Q6, Q7 or Q9 and B1, B2, B3 and B4 are carbon atoms at the

same time, then Ym is other than 3-chloro-2-methyl group, 3-chloro-2,6-diethyl group, 5-chloro-2-methyl group, 2,6-diethyl group, 4-chloro-2-fluoro group and 2-ethyl-6-methyl group,

- (2) when Het is Q4 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 2,5-dichloro group, 2,4-diffuoro group, 2,6-diffuoro group, 3-chloro-2-methyl group, 5-chloro-2-methyl group, 5-fluoro-2-methyl group, 2,6-dimethyl group, 2,6-diethyl group, 2-ethyl-6-methyl group, 2-methoxy-5-nitro group, 2-methoxy-5-methyl group, 2,6-diethoxy group, 3-bromo-2-methyl group, 3-fluoro-2-methyl group, 3-iodo-2-methyl group, 3-cyano-2-methyl group, 3-diffuoromethoxy-2-methyl group, 5-chloro-2-ethyl group, 2,5-dimethyl group, 2,3-dichloro group, 3-chloro-2,6-diethyl group, 4-trifluoromethyl group, 3-methoxycarbonyl-2-methyl group, 3-trifluoromethyl-2-methyl group, 3,5-dichloro-2,6-diethyl group, 3,4-dichloro group, 3-(methoxycarbonylmethyloxy)-2-methyl group, 2-methyl-3-nitro group and 4-trifluoromethoxy group,
- (3) when Het is Q9, R² and R³ are hydrogen atoms at the same time, Xn is a 2-phenyl group, R¹ is a n-propyl group or an i-propyl group and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 4-pentafluoroethyl-2-methyl group, and
- (4) when Het is Q10 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 5-chloro-2-methyl group, 5-fluoro-2-methyl group [and] , 2,5-dimethyl group and 2,6-diethyl group, and
- (5) when Het is Q10 and B¹, B², B³ and B⁴ are carbon atoms at the same time, Xn is other than 5,6-dimethyl group }.

IN THE UNITED STATES PATENT AND TRADEMARK OFFICE



In re PATENT APPLICATION of KATSUHIRA, et al.

Group Art Unit: Not Assigned

Appln. No.: Not Assigned

Examiner: Not Assigned

Filed: December 3, 2001

International Appln. No. PCT/JP00/04136

Title: HETEROCYCLIC DICARBOXYLIC ACID DIAMIDE DERIVATIVES,

AGRICULTURAL AND HORTICULTURAL INSECTICIDES AND

METHOD OF USING THE SAME

December 19, 2001

FIRST PRELIMINARY AMENDMENT: TO BE ENTERED PRIOR TO CALCULATION OF FILING FEE

Hon. Commissioner of Patents and Trademarks Washington, D.C. 20231

Sir:

Please enter the following Preliminary Amendment of the subject new application prior to calculation of the fee for filing the application.

IN THE CLAIMS:

Please amend claim 6 as follows (see the attached Appendix for the changes made to effect the below claim):

Claim 6. (Amended) An agricultural and horticultural insecticide characterized by containing a heterocyclic dicarboxylic acid diamide derivative according to claim 1 as an active ingredient.

REMARKS

This Preliminary Amendment revises the multiple dependent claims to be single dependent claims and thus reduce the filing fee for the subject application. No new matter has been added.

Entry of this amendment and favorable consideration of this application are respectfully requested.

Respectfully submitted,

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APPENDIX SHOWING REVISIONS OF CLAIMS

Proposed Amendments To Claim 6 Showing Deletions And Insertions.

Claim 6. (Amended) An agricultural and horticultural insecticide characterized by containing a heterocyclic dicarboxylic acid diamide derivative according to [any one of claims 1 to 5] <u>claim 1</u> as an active ingredient.

581 Recard 19 DEC 2001

DESCRIPTION

HETEROCYCLIC DICARBOXYLIC ACID DIAMIDE DERIVATIVES, AGRICULTURAL AND HORTICULTURAL INSECTICIDES AND METHOD OF USING THE SAME

TECHNICAL FIELD

The present invention relates to heterocyclic dicarboxylic acid diamide derivatives, agricultural and horticultural insecticides containing any of said compounds as an active ingredient, and their usage.

BACKGROUND ART

JP-A-6-25190, JP-A-9-323974, WO9944992, JP-A-12-7661, JP-A-12-103708 and the like disclose compounds constituting a portion of the heterocyclic dicarboxylic acid diamide derivative of the present invention but do not describe or suggest the usefulness of the disclosed compounds as agricultural and horticultural insecticides at all.

The present inventors earnestly investigated 15 in order to develop a novel agricultural and horticultural insecticide, and consequently found that heterocyclic dicarboxylic acid diamide derivatives represented by the general formula (I) of the present invention are novel compounds not known in any literature and that not only these compounds but also the compounds disclosed in the prior arts can be put to a novel use, whereby the present invention has been

accomplished.

DISCLOSURE OF THE INVENTION

The present invention relates to heterocyclic dicarboxylic acid diamide derivatives represented by the general formula (I):

{wherein R^1 , R^2 and R^3 , which may be the same or different, are hydrogen atoms, (C3-C6)cycloalkyl groups, halo (C_3-C_6) cycloalkyl groups or $-A^1-(R^4)$ r (wherein A^1 is a (C_1-C_8) alkylene group, a (C_3-C_6) alkenylene group or a 10 (C_3-C_6) alkynylene group, R^4 , which may be the same or different, are hydrogen atoms; halogen atoms; cyano groups; nitro groups; halo(C1-C6)alkyl groups; (C3- C_6) cycloalkyl groups; halo (C_3-C_6) cycloalkyl groups; (C_1-C_6) C_6) alkoxycarbonyl groups; di (C_1-C_6) alkoxyphosphoryl 15 groups whose (C_1-C_6) alkoxy groups may be the same or different; $di(C_1-C_6)$ alkoxythiophosphoryl groups whose (C_1-C_6) alkoxy groups may be the same or different; diphenylphosphino groups; diphenylphosphono groups; phenyl groups; substituted phenyl groups having one or 20 more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl

groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl

- groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups,
- halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^2-R^5$ (wherein A^2 is -O-, -S-, -SO-, $-SO_2-$, $-N(R^6)-$ (wherein R^6
- is a hydrogen atom; a (C_1-C_6) alkylcarbonyl group; a halo (C_1-C_6) alkylcarbonyl group; a (C_1-C_6) alkoxycarbonyl group; a phenylcarbonyl group; a substituted phenylcarbonyl group having one or more substituents which may be the same or different and are selected
- from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6)
- C₆) alkylsulfonyl groups; a phenyl(C_1 - C_4) alkoxycarbonyl group; a substituted phenyl(C_1 - C_4) alkoxycarbonyl group having on the ring one or more substituents which may be the same or different and are selected from halogen

atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a (C_1-C_6) alkylsulfonyl group; or a halo (C_1-C_6) alkylsulfonyl group), -C(=O) or $-C(=NOR^7)$ (wherein R^7 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo (C_1-C_6) a halo

- 10 C_6) alkenyl group; a (C_3-C_6) alkynyl group; a cyclo (C_3-C_6) alkyl group; a phenyl (C_1-C_4) alkyl group; or a substituted phenyl (C_1-C_4) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_4) alkyl group having on the ring
- 15 C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl
- groups), and R^5 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo (C_1-C_6) alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) alkenyl group; a (C_3-C_6) alkynyl group; a halo (C_3-C_6) alkynyl group; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxy (C_1-C_6) alkyl
- group; a (C_1-C_6) alkylthio (C_1-C_6) alkyl group; a formyl group; a (C_1-C_6) alkylcarbonyl group; a halo (C_1-C_6) alkylcarbonyl group; a (C_1-C_6) alkoxycarbonyl group; a mono (C_1-C_6) alkylaminocarbonyl group; a di (C_1-C_6)

 C_6) alkylaminocarbonyl group whose (C_1-C_6) alkyl groups may be the same or different; a mono (C_1-C_6) alkylaminothiocarbonyl group; a di (C_1-C_6) alkylaminothiocarbonyl group whose (C_1-C_6) alkyl groups may be the same or different;

- a di (C_1-C_6) alkoxyphosphoryl group whose (C_1-C_6) alkoxy groups may be the same or different; a di (C_1-C_6) alkoxythiophosphoryl group whose (C_1-C_6) alkoxy groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substi-
- tuents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups,
- halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenyl (C_1-C_4) alkyl group; a substituted phenyl (C_1-C_4) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen
- atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl
- groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl

groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and r is an integer of 1 to 4),

provided that $\ensuremath{R^1}$ and $\ensuremath{R^2}$ are not hydrogen atoms at the same time,

R¹ and R² may form a 4 to 7 membered ring by

10 combining to each other, in which the ring may contain
the same or different 1 to 3 hetero atoms selected from
the group consisting of oxygen atom, sulfur atom and
nitrogen atom,

Het is a heterocyclic group represented by any of the following formulas Q1 to Q22:

(wherein X, which may be the same or different, are halogen atoms; cyano groups; nitro groups; (C_3-C_6) cycloalkyl groups; halo (C_3-C_6) cycloalkyl groups; tri (C_1-C_6) alkylsilyl groups whose (C_1-C_6) alkyl groups may be the same or different; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6)

- 10 C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the
- same or different and are selected from halogen atoms, $(C_1-C_6) \, \text{alkyl groups, halo} \, (C_1-C_6) \, \text{alkyl groups, } \, (C_1-C_6) \, \text{alkoxy groups, halo} \, (C_1-C_6) \, \text{alkoxy groups, } \, (C_1-C_6) \, \text{alkylthio groups, halo} \, (C_1-C_6) \, \text{alkylthio groups, halo} \, (C_1-C_6) \, \text{alkylsulfinyl groups, halo} \, (C_1-C_6)$
- 20 (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ [wherein A^3 is -O-, -S-, -SO-, $-SO_2-$, $-N(R^6)-$ (wherein R^6 is as defined above), -C(=O)-, $-C(=NOR^7)-$ (wherein R^7 is as defined above), a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6)
- C₆) alkenylene group, a halo (C_2-C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R⁸ is as follows:
 - (1) when A^3 is -O-, -S-, -SO-, $-SO_2-$ or $-N(R^6)-$ (wherein

 R^6 is as defined above), then R^8 is a halo(C_3 - C_6) cycloalkyl group; a halo(C_3 - C_6) cycloalkenyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1 - C_6) alkyl groups, halo(C_1 - C_6) alkyl groups, (C_1 - C_6) alkoxy groups, halo(C_1 - C_6) alkoxy groups, (C_1 - C_6) alkylthio groups, halo(C_1 - C_6) alkylthio groups, (C_1 - C_6) alkylsulfinyl groups, halo(C_1 - C_6) alkylsulfinyl groups, (C_1 - C_6) alkylsulfinyl groups,

- groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy
- groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^4-R^9$ (wherein A^4 is a (C_1-C_6) alkylene group,
- a halo (C_1-C_6) alkylene group, a (C_3-C_6) alkenylene group, a halo (C_3-C_6) alkenylene group, a (C_3-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R^9 is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxycarbonyl group; a
- 25 phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups,

halo(C₁-C₆) alkoxy groups, (C₁-C₆) alkylthio groups,
halo(C₁-C₆) alkylthio groups, (C₁-C₆) alkylsulfinyl groups,
halo(C₁-C₆) alkylsulfinyl groups, (C₁-C₆) alkylsulfonyl
groups and halo(C₁-C₆) alkylsulfonyl groups; or -A⁵-R¹⁰

(wherein A⁵ is -O-, -S-, -SO-, -SO₂- or -C(=O), and R¹⁰
is a (C₁-C₆) alkyl group; a halo(C₁-C₆) alkyl group; a (C₃C₆) alkenyl group; a halo(C₃-C₆) alkenyl group; a (C₃C₆) cycloalkyl group; a halo(C₃-C₆) cycloalkyl group; a
phenyl group; a substituted phenyl group having one or
more substituents which may be the same or different
and are selected from halogen atoms, (C₁-C₆) alkyl
groups, halo(C₁-C₆) alkyl groups, (C₁-C₆) alkoxy groups,
halo(C₁-C₆) alkoxy groups, (C₁-C₆) alkylsulfinyl groups,

- halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkylsulfonyl groups; a
- 20 C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl
- 25 groups)),
 - (2) when A^3 is -C(=0) or $-C(=NOR^7)$ (wherein R^7 is as defined above), then R^8 is a hydrogen atom; a (C_1 - C_6) alkyl group; a halo (C_1 - C_6) alkyl group; a (C_2 -

- C_6) alkenyl group; a halo (C_2-C_6) alkenyl group; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxy group; a (C_1-C_6) alkylamino group; a di (C_1-C_6) alkylamino group whose
- 5 (C_1-C_6) alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups,
- halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenylamino group; a substituted phenylamino group
- having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6)
- C₆) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected
- from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl

groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups, and

- (3) when A^3 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6)
- 5 C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, then R^8 is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxycarbonyl group; a tri (C_1-C_6) alkylsilyl group whose (C_1-C_6) alkyl groups may
- be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkoxy groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6)
- C₆) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents
- which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl
- groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^6-R^{11}$ (wherein A^6 is -O-, -S-, -SO- or $-SO_2-$, and R^{11} is a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a phenyl group; a

substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6)

- 5 C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group
- having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl
- groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^7-R^{12}$ (wherein A^7 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6) alkenylene group, a (C_2-C_6) alkynylene group or
- a halo (C_3-C_6) alkynylene group, and R^{12} is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxy group; a halo (C_1-C_6) alkoxy group; a (C_1-C_6) alkylthio group; a halo (C_1-C_6) alkylthio group; a (C_1-C_6) alkylthio group;
- a halo (C_1-C_6) alkylsulfinyl group; a (C_1-C_6) alkylsulfonyl group; a halo (C_1-C_6) alkylsulfonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are

selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups,

- halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenoxy group; a substituted phenoxy group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups,
- halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenylthio
- group; a substituted phenylthio group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups,
- 20 halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same
- or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl

groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups)), and n is an integer of 0 to 3,

X may form a condensed ring by combining together with the adjacent atoms in the heterocyclic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups;

- 10 C_6) alkoxy groups; (C_1-C_6) alkylthio groups; halo (C_1-C_6) alkylthio groups; (C_1-C_6) alkylsulfinyl groups; halo (C_1-C_6) alkylsulfinyl groups; (C_1-C_6) alkylsulfonyl groups; halo (C_1-C_6) alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more
- substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups,
- halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl
- groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl

groups and halo (C_1-C_6) alkylsulfonyl groups,

W is O, S or N-R¹³ (wherein R¹³ is a (C₁-C₆) alkyl group; a halo (C₁-C₆) alkyl group; a (C₃-C₆) alkenyl group; a halo (C₃-C₆) alkenyl group; a (C₁-C₆) alkynyl group; a halo (C₃-C₆) alkynyl group; a (C₁-C₆) alkoxy group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C₁-C₆) alkyl groups, halo (C₁-C₆) alkyl groups, (C₁-

- 10 C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenyl (C_1-C_6) alkyl group; or a substituted
- phenyl(C_1-C_6) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1-C_6) alkylthio groups,
- C₆) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and p and q, which may be the same or different, are integers of 0 to 1),
- B^1 , B^2 , B^3 and B^4 , which may be the same or different, are carbon atoms or nitrogen atoms,

Y, which may be the same or different, are halogen atoms; cyano groups; nitro groups; halo(C_3 -

 C_6) cycloalkyl groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6)

- 5 C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic
- groups having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1-C_6) \, \text{alkyl groups, halo} \, (C_1-C_6) \, \text{alkyl groups, } \, (C_1-C_6) \, \text{alkoxy groups, halo} \, (C_1-C_6) \, \text{alkoxy groups, halo} \, (C_1-C_6) \, \text{alkylthio groups$
- C₆) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ (wherein A^3 and R^8 are as defined above), and m is an integer of 1 to 5,

Y may form a condensed ring by combining

20 together with the adjacent carbon atoms in the aromatic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C₁-C₆) alkyl groups; halo(C₁-C₆) alkyl groups; (C₁-C₆) alkoxy groups; halo(C₁-C₆) alkoxy groups; (C₁-C₆) alkylthio groups; (C₁-C₆) alkylthio groups; (C₁-C₆) alkylsulfinyl groups;

 C_6) alkoxy groups; (C_1-C_6) alkylthio groups; halo (C_1-C_6) alkylthio groups; (C_1-C_6) alkylsulfinyl groups; halo (C_1-C_6) alkylsulfinyl groups; (C_1-C_6) alkylsulfonyl groups; halo (C_1-C_6) alkylsulfonyl groups; phenyl group;

substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6)

- 5 C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one
- or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthio groups,
- halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups, and each of Z^1 and Z^2 is an oxygen atom or a sulfur atom,

provided that:

- (1) when Het is Q2, Q6, Q7 or Q9 and B¹, B², B³ and B⁴
 20 are carbon atoms at the same time, then Ym is other than 3-chloro-2-methyl group, 3-chloro-2,6-diethyl group, 5-chloro-2-methyl group, 2,6-diethyl group, 4-chloro-2-fluoro group and 2-ethyl-6-methyl group,
 (2) when Het is Q4 and B¹, B², B³ and B⁴ are carbon
- atoms at the same time, then Ym is other than 2,5-dichloro group, 2,4-difluoro group, 2,6-difluoro group, 3-chloro-2-methyl group, 5-chloro-2-methyl group, 5-fluoro-2-methyl group, 2,6-diethyl

group, 2-ethyl-6-methyl group, 2-methoxy-5-nitro group, 2-methoxy-5-methyl group, 2,6-diethoxy group, 3-bromo-2-methyl group, 3-fluoro-2-methyl group, 3-iodo-2-methyl group, 3-cyano-2-methyl group, 3-

- difluoromethoxy-2-methyl group, 5-chloro-2-ethyl group, 2,5-dimethyl group, 2,3-dichloro group, 3-chloro-2,6-diethyl group, 4-trifluoromethyl group, 3-methoxycarbonyl-2-methyl group, 3-trifluoromethyl-2-methyl group, 3,5-dichloro-2,6-diethyl group, 3,4-
- dichloro group, 3-(methoxycarbonylmethyloxy)-2-methyl group, 2-methyl-3-nitro group and 4-trifluoromethoxy group,
 - (3) when Het is Q9, R^2 and R^3 are hydrogen atoms at the same time, Xn is a 2-phenyl group, R^1 is a n-propyl
- 15 group or an i-propyl group and B^1 , B^2 , B^3 and B^4 are carbon atoms at the same time, then Ym is other than 4-pentafluoroethyl-2-methyl group, and
 - (4) when Het is Q10 and B^1 , B^2 , B^3 and B^4 are carbon atoms at the same time, then Ym is other than 5-chloro-
- 20 2-methyl group, 5-fluoro-2-methyl group and 2,5-dimethyl group}, agricultural and horticultural insecticides and their usage.

MODE FOR CARRYING OUT THE INVENTION

In the definition of the general formula (I)

25 for the heterocyclic dicarboxylic acid diamide derivative of the present invention, the term "halogen atom" means a chlorine atom, a bromine atom, an iodine

atom or a fluorine atom. The term " (C_1-C_6) alkyl" means a linear or branched alkyl group of 1 to 6 carbon atoms, such as methyl, ethyl, n-propyl, i-propyl, n-butyl, i-butyl, s-butyl, t-butyl, n-pentyl, n-hexyl or the like.

- The term "halo(C_1 - C_6) alkyl" means a substituted linear or branched alkyl group of 1 to 6 carbon atoms having as the substituent(s) one or more halogen atoms which may be the same or different. The term "halo(C_3 - C_6) cycloalkyl" means a substituted alicyclic hydrocarbon
- group of 3 to 6 carbon atoms having as the substituent(s) one or more halogen atoms which may be the same or different. The term " (C_1-C_8) alkylene" means a linear or branched alkylene group of 1 to 8 carbon atoms, such as methylene, ethylene, propylene,
- 15 trimethylene, dimethylmethylene, tetramethylene, isobutylene, dimethylethylene, octamethylene or the like. The term "heterocyclic group" means a heterocyclic group such as pyridyl group, pyridine-Noxide group, pyrimidinyl group, furyl group,
- tetrahydrofuryl group, thienyl group, tetrahydrothienyl group, tetrahydropyranyl group, tetrahydrothiopyranyl group, oxazolyl group, isoxazolyl group, oxadiazolyl group, thiazolyl group, isothiazolyl group, thiadiazolyl group, imidazolyl group, triazolyl group,
- 25 pyrazolyl group or the like. As to the passage " R^1 and R^2 may form a 4 to 7 membered ring by combining to each other, in which the ring may contain the same or different 1 to 3 hetero atoms selected from the group

consisting of oxygen atom, sulfur atom and nitrogen atom", the 4 to 7 membered ring includes, for example, azetidine ring, pyrrolidine ring, pyrroline ring, piperidine ring, imidazolidine ring, imidazoline ring, oxazolidine ring, thiazolidine ring, isoxazolidine ring, isothiazolidine ring, tetrahydropyridine ring, piperazine ring, morpholine ring, thiomorpholine ring, dioxazine ring, dithiazine ring, indole ring, benzo[b]furan ring, benzo[b]thiophene ring, quinoline ring, isoquinoline ring, naphthyridine ring and quinoxaline ring.

As to the passage "X may form a condensed ring by combining together with the adjacent atoms in the heterocyclic ring", the condensed ring refers to, for example, indole ring, benzo[b] furan ring, benzo[b] thiophene ring, quinoline ring, isoquinoline ring, naphthyridine ring, quinoxaline ring or cinnoline ring.

As to the passage "Y may form a condensed ring by combining together with the adjacent carbon atoms in the aromatic ring", the condensed ring refers to a condensed ring such as naphthalene, tetrahydronaphthalene, indene, indan, quinoline, quinazoline, chroman, isochroman, indole, indoline, benzodioxane,

25 benzodioxole, benzofuran, dihydrobenzofuran, benzothiophene, dihydrobenzothiophene, benzooxazole, benzothiazole, benzoimidazole, indazole or the like.

The heterocyclic dicarboxylic acid diamide

derivative of the general formula (I) of the present invention contains an asymmetric carbon atom or an asymmetric center in the structural formula in some cases, and has optical isomers and diastereomers in some cases. The present invention includes all of the individual optical isomers and mixtures containing the optical isomers in any proportions. The heterocyclic dicarboxylic acid diamide derivative of the general formula (I) of the present invention has geometrical isomers due to a carbon-carbon double bond or a carbon-nitrogen double bond in the structural formula in some cases. The present invention includes all of the individual geometrical isomers and mixtures containing the geometrical isomers in any proportions.

dicarboxylic acid diamide derivative of the general formula (I) of the present invention are a group of compounds in which R¹ is a (C₁-C₀)alkyl group, a (C₁-C₀)alkylthio(C₁-C₀)alkyl group, a (C₁-C₀)alkylthio(C₁-C₀)alkyl group, a (C₁-C₀)alkyl-sulfonyl(C₁-C₀)alkyl group; R² and R³, which may be the same or different, are hydrogen atoms or methyl groups; Het is a pyridine ring represented by Q1, Q2, Q3 or Q4; X, which may be the same or different, are halogen atoms, nitro groups, halo(C₁-C₀)alkyl groups, halo(C₁-C₀)alkyl groups; n is an integer of 0 to 2; p is an integer of 0 or 1; all of B¹, B², B³ and B⁴ are carbon atoms, or all of B¹, B² and B⁴ are carbon atoms

and B³ is a nitrogen atom; Y, which may be the same or different, are halogen atoms, (C₁-C₆) alkyl groups, halo(C₁-C₆) alkyl groups, halo(C₁-C₆) alkyl groups, halo(C₁-C₆) alkylthio groups or halo(C₁-C₆) alkoxyhalo(C₁-C₆) alkoxy groups; as to the substitution position of Y and the number m of substituents Y, the substitution is di-substitution at the 2- and 3-positions or the 2- and 4-positions in relation to the binding position of the amide group, or tri-substitution at the 2-, 3- and 4- positions or the 2-, 4- and 5-positions in relation to the binding position of the amide group; and each of Z¹ and Z² is an oxygen atom.

A group of more preferable compounds are compounds in which R1 is an i-propyl group, a t-butyl 15 group, a methylthio (C_3-C_4) alkyl group, a methyl $sulfinyl(C_3-C_4)$ alkyl group or a methylsulfonyl($C_3 C_4$) alkyl group; each of R^2 and R^3 is a hydrogen atom; Het is a pyridine ring represented by Q1, Q2, Q3 or Q4; X is halogen atom; n is an integer of 0 to 1; p is an integer of 0 or 1; all of B^1 , B^2 , B^3 and B^4 are carbon 20 atoms; Y, which may be the same or different, are chlorine atoms, methyl groups, trifluoromethyl groups, pentafluoroethyl groups, heptafluoropropyl groups, heptafluoroisopropyl groups, trifluoromethoxy groups or 1-trifluoromethyl-2,2,2-trifluoroethoxy groups; as to the substitution position of Y and the number m of substituents Y, the substitution is di-substitution at the 2- and 4-positions in relation to the binding

position of the amide group; and each of Z^1 and Z^2 is an oxygen atom.

The heterocyclic dicarboxylic acid diamide derivative of the general formula (I) of the present invention can be produced, for example, by any of the production processes schematically shown below.

Production process 1.

$$\begin{array}{c} \text{Xn} \\ \text{CO}_2 \text{R} \\ \text{(VIII)} \\ \\ \text{(VIII)} \\ \\ \text{(VIII)} \\ \\ \text{(VIII)} \\ \\ \text{(VII)} \\ \\ \text{(IV-2)} \\ \\ \text{(IV-2)} \\ \\ \text{(IV-2)} \\ \\ \text{(IV-2)} \\ \\ \text{(IIII)} \\ \\ \text{(III)} \\ \\ \text{(III)} \\ \\ \text{(IIII)} \\ \\ \text{(III)} \\ \\ \text{(IIII)} \\ \\ \text{(IIIII$$

the imide.

wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, m and n are as defined above, hal is a halogen atom, and R is a (C_1-C_3) alkyl group.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) can be produced by hydrolyzing a diester of the general formula (VIII) in the presence of an acid or an alkali to obtain a dicarboxylic acid of the general formula (VII), converting said dicarboxylic acid to an acid anhydride 10 (VI) in the presence of a dehydrating agent, reacting the acid anhydride (VI) with a substituted aromatic amine of the general formula (V) in the presence or absence of an inert solvent to obtain amides of the general formulas (IV-1) and (IV-2), reacting said amides with a dehydrating agent in the presence or absence of an inert solvent after or without isolating the amides, to obtain an imide of the general formula (III), and then reacting said imide with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, in the presence or 20 absence of an inert solvent after or without isolating

 solvents.

As the base used for the hydrolysis,
hydroxides of alkali metals, such as sodium hydroxide,
potassium hydroxide and the like can be used. As to

5 the amount of the base used, the base may be used in an
amount properly chosen in the range of 2 to 10
equivalents per equivalent of the diester of the
general formula (VIII). As the acid, there can be
used, for example, inorganic acids such as hydrochloric

10 acid, sulfuric acid, etc.; and organic acids such as
trifluoroacetic acid, etc. The amount of the acid used
may be a catalytic amount relative to the diester of
the general formula (VIII) and ranges from 0.001 to 0.1
equivalent per equivalent of the diester.

As to the reaction temperature, the reaction can be carried out at room temperature or while refluxing the inert solvent used. Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, the reaction may be carried out for a time properly chosen in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by a recrystallization method, a distillation method, a column chromatographic method, etc., whereby the desired compound can be produced.

It is also possible to subject the desired compound to the subsequent reaction without isolation after completion of the reaction.

As an inert solvent usable in this reaction, any inert solvent may be used so long as it does not markedly inhibit the progress of the reaction. There can be used, for example, halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride, etc.; aromatic hydrocarbons such as benzene, toluene, xylene, chlorobenzene, etc.; acyclic or cyclic ethers such as Methyl Cellosolves, diethyl ether, diisopropyl ether, dioxane, tetrahydrofuran, etc.; and organic acids such as acetic acid, trifluoroacetic acid, etc.

These inert solvents may be used singly or as a mixture thereof.

The dehydrating agent may be used in excess to serve also as the inert solvent.

As the dehydrating agent, there can be used

dehydrating agents such as acetic anhydride,

trifluoroacetic anhydride, etc. As to the amount of

these dehydrating agents used, they may be used in an

amount properly chosen in the range of 1 mole to excess

moles per mole of the compound of the general formula

(VII), and they are preferably used in an amount

equimolar with the amount of the compound of the

general formula (VII).

15

20

The reaction temperature may be properly chosen in the range of room temperature to the boiling point of the inert solvent used. When no inert solvent is used, it is sufficient that the reaction is carried out below the boiling point of the dehydrating agent used.

Although the reaction time is varied depending on the reaction temperature, the scale of reaction, and the like, the reaction may be carried out in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by a recrystallization method, a distillation method, a column chromatographic method, etc., whereby the desired

It is also possible to subject the desired compound to the subsequent reaction without isolation after completion of the reaction.

compound can be produced.

(1-3). General formula (VI) \rightarrow general formula (IV-1) + general formula (IV-2)

As the inert solvent usable in this reaction, any inert solvent may be used so long as it does not markedly inhibit the progress of the reaction. There can be exemplified inert solvent including aromatic hydrocarbons such as benzene, toluene, xylene, etc.;

halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride, chlorobenzene, dichlorobenzene, etc.; acyclic or cyclic ethers such as diethyl ether, dioxane, tetrahydrofuran, etc.; esters such as ethyl acetate, etc.; amides such as dimethylformamide, dimethylacetamide, etc.; acids such as acetic acid, etc.; dimethyl sulfoxide; 1,3-dimethyl-2-imidazolidinone; and water. These inert solvents may be used singly or as a mixture of two or more thereof.

Since the reaction is an equimolar reaction, it is sufficient that the reactants are used in equimolar amounts, though either of them may be used in excess. If necessary, the reaction may be carried out under dehydrating conditions.

As to the reaction temperature, the reaction can be carried out at room temperature or while refluxing the inert solvent used. Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, the reaction may be carried out for a time properly chosen in the range of several minutes to 48 hours.

After completion of the reaction, the desired compounds are isolated from the reaction system containing the desired compounds by a conventional method, and if necessary, purified by a recrystallization method, a distillation method, a column chromatographic method, etc., whereby the desired compounds can be produced.

It is also possible to subject the desired compounds to the subsequent reaction without isolation after completion of the reaction.

(1-4). General formula (IV-1) + general formula (IV-2) \rightarrow general formula (III)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in (1-2).

It is also possible to subject the desired compounds to the subsequent reaction without isolation after completion of the reaction.

(1-5). General formula (III) \rightarrow general formula (I-1)

As the inert solvent usable in this reaction,
pyridines can be used besides, for example, the inert
solvents exemplified in (1-2).

Since the reaction is an equimolar reaction, it is sufficient that the amine of the general formula (II-1) or the amine salt of the general formula (II-2) is used in an amount equimolar with the amount of the imide of the general formula (III), though the amine or the amine salt may be used in excess.

When the amine salt of the general formula

(II-2) is used in the reaction, a base is necessary in

order to produce free amine in the reaction system. As

the base, an inorganic base or an organic base can be

used. The inorganic base includes, for example,

hydroxides and carbonates of alkali metals, such as sodium hydroxide, potassium hydroxide, sodium carbonate and potassium carbonate. The organic base includes, for example, triethylamine, pyridine, 4-(dimethyl-amino)pyridine and 1,8-diazabicyclo[5.4.0]-7-undecene. As to the amount of these bases used, they may be used in an amount properly chosen in the range of 1 mole to excess moles per mole of the amine salt of the general formula (II-2).

The reaction temperature may be properly chosen in the range of -10°C to the boiling point of the inert solvent used, and the reaction is preferably carried out in the range of 0°C to 150°C .

Although the reaction time is varied

15 depending on the reaction temperature, the scale of reaction, and the like, the reaction may be carried out in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system

20 containing the desired compound by a conventional method, and if necessary, purified by a recrystal-lization method, a distillation method, a column-chromatographic method, etc., whereby the desired compound can be produced.

The compound of the general formula (VIII), i.e., the starting material in the reaction can be produced according to a well-known process, for example, the processes described in, for instance, J.

- Am. Chem. Soc., <u>63</u>, 1762 (1941), J. Heterocyclic Chem., <u>21</u>, 1431 (1984), J. Indian Chem. Soc., <u>1982</u>, 1372, J. Org. Chem., <u>14</u>, 723 (1949), Heterocycles, <u>27</u>, 1489 (1988), J. Am. Chem. Soc., <u>78</u>, 2220 (1956), J. Prakt.
- 5 Chem., 311, 807 (1969), Tetrahedron, 36, 1801 (1980), JP-A-6-122684, U.S. Patent 3,414,580, ditto 3,686,171, J. Med. Chem., 27, 1396 (1984), J. Heterocyclic Chem., 12, 1303 (1975), ditto 15, 1477 (1978), ditto 16, 1141 (1979), ditto 17, 443 (1982), ditto 21, 689 (1984),
- 10 Beil., <u>25III</u>, 2028, JP-A-52-77086, J. Am. Chem. Soc., <u>81</u>, 2456 (1956), J. Org. Chem., <u>37</u>, 3224 (1972), JP-A-62-175480, JP-A-62-230782, JP-A-60-69083, JP-A-60-185783, JP-A-61-109790, JP-A-62-277385, JP-A-63-295575, JP-A-63-99067, JP-A-64-75474, JP-A-64-90118, Yakugaku
- 15 Zasshi, 84, 416 (1964), Chem. and Pharm. Bull., 5, 277
 (1957), Chem. Research (S), 1989, 196, Chem. Pharm.
 Bull., 20, (7), 1513 (1972), J. Heterocyclic Chem., 27,
 579 (1990), Tetrahedron, 53 (42), 14497 (1997), ditto
 41 (7), 1199 (1985), Chem. Ber., 107, 3036 (1974), J.
- 20 Heterocyclic Chem., <u>23</u>, 1103 (1986), ditto <u>5</u>, 125 (1968), J. Org. Chem., <u>26</u>, 468 (1961), etc.

10

Production process 2.

X'n

$$B^1 = B^2$$
 Ym

 $B^1 = B^2$ Ym

 $B^1 = B^2$ Ym

(III-1)

 $R^1 R^2 NH$ (II-1)

 $R^1 R^2 NH \bullet Hhal$ (II-2)

 $R^1 R^2 NH \bullet Hhal$ (II-2)

wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and n are as defined above, and X' is a halogen atom or a nitro group, provided that X is other than hydrogen atom and nitro group.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) can be produced by reacting a heterocyclic dicarboxylic acid imide derivative of the general formula (III-1) with a reactant corresponding to X, in the presence of an inert solvent to obtain a heterocyclic dicarboxylic acid imide derivative of the general formula (III), and reacting said heterocyclic dicarboxylic acid imide derivative (III) with an amine or its salt of the general formula (III-1),

respectively, after or without isolating the derivative (III).

- 10 (2-2). General formula (III) \rightarrow general formula (I-1) In the case of this reaction, the production can be carried out according to production process (1-5).

Production process 3.

$$\begin{array}{c} B^1 = B^2 \text{ Ym} \\ O_2N \\ \text{Het} \\ O_2N \\ \text{CO}_2H \\ \text{CO}_2$$

wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and n are as defined above.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) can be produced by reacting a heterocyclic dicarboxylic acid anhydride derivative of the general formula (VI-1) with an aromatic amine of the general formula (V) in the presence of an inert solvent to obtain amides of the

general formulas (IV-1') and (IV-2'), reacting said amides with a dehydrating agent in the presence or absence of an inert solvent after or without isolating the amides, to obtain a heterocyclic dicarboxylic acid

- imide derivative of the general formula (III-2), subjecting said heterocyclic dicarboxylic acid imide derivative (III-2) to catalytic reduction with hydrogen after or without isolation to obtain a heterocyclic dicarboxylic acid imide derivative of the general
- formula (III-3), diazotizing said heterocyclic dicarboxylic acid imide derivative (III-3) after or without isolation, adding a metal salt thereto to obtain a heterocyclic dicarboxylic acid imide derivative of the general formula (III), and then
- reacting said heterocyclic dicarboxylic acid imide derivative (III) with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, after or without isolating the derivative (III).
- 20 (3-1). General formula (VI-1) \rightarrow general formula (IV-1') + general formula (IV-2')

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-3).

25 (3-2). General formula (IV-1') + general formula (IV-2') \rightarrow general formula (III-2)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-4).

(3-3). General formula (III-2) \rightarrow general formula (III-5)

As an inert solvent usable in this reaction, any inert solvent may be used so long as it does not markedly inhibit the progress of the reaction. There can be exemplified alcohols such as methanol, ethanol, propanol, etc.; acyclic or cyclic ethers such as diethyl ether, dioxane, tetrahydrofuran, etc.; and acids such as acetic acid, etc. These inert solvents may be used singly or as a mixture of two or more thereof.

As a catalyst for the catalytic reduction used in the reaction, there can be exemplified palladium-carbon, Raney nickel, palladium black and platinum black. As to the amount of the catalyst used, the catalyst may be used in an amount properly chosen in the range of 0.1 to 10 wt% based on the weight of the heterocyclic dicarboxylic acid imide derivative of the general formula (III-2). The reaction is carried out under a hydrogen atmosphere, and the reaction may be carried out at a hydrogen pressure properly chosen in the range of 1 to 10 atmospheric pressure.

As to the reaction temperature, the reaction may be carried out at room temperature to the reflux

temperature of the inert solvent used. Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, it may be properly chosen in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystal—

10 lization, column chromatography, etc., whereby the desired compound can be produced. It is also possible to subject the desired compound to the subsequent reaction without isolation from the reaction system.

- As an inert solvent used in this reaction, acidic solvents can be used. There can be exemplified an aqueous hydrochloric acid solution, an aqueous hydrobromic acid solution, an aqueous hydroiodic acid solution, an aqueous solution, an aqueous sulfuric acid solution, acetic acid and trifluoroacetic acid. These acidic solvents may be used singly or as a mixture of two or more thereof. In addition, these acidic solvents may be used in admixture with ethers such as tetrahydrofuran, dioxane, etc.
- As a diazotizing agent, there can be exemplified diazotizing agents such as sodium nitrite, nitrosyl hydrogensulfate, alkyl nitrites, etc. As to

the amount of these diazotizing agents used, the reaction may be carried out by properly choosing the amount in the range of an amount equal to that of the heterocyclic dicarboxylic acid imide derivative of the general formula (III-3) to an excess amount over this derivative.

As to the reaction temperature, the reaction may be carried out at $-50^{\circ}\mathrm{C}$ to room temperature or at the reflux temperature of the inert solvent used.

10 Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, it may be properly chosen in the range of several minutes to 48 hours.

As the metal salt added after the production of the diazonium salt, there can be used metal salts such as cuprous chloride, cuprous bromide, potassium iodide, copper cyanide, potassium xanthate, mercaptan sodium, etc. As to the amount of the metal salt used, the reaction may be carried out by properly choosing

the amount in the range of 1 equivalent to excess equivalents per equivalent of the heterocyclic dicarboxylic acid imide derivative of the general formula (III-3).

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystallization, column chromatography, etc., whereby the

desired compound can be produced. It is also possible to subject the desired compound to the subsequent reaction without isolation from the reaction system.

In the case of the reaction, the production can be carried out according to the process described in Org. Synth., IV, 160 (1963), ditto III, 809 (1959), J. Am. Chem. Soc., 92, 3520 (1970), etc.

(3-5). General formula (III) \rightarrow general formula (I-1)

In the case of this reaction, the desired

compound can be produced in a manner similar to that described in production process (1-5).

Production process 4.

NO2
Heth
$$B^1 = B^2$$
 Ym
$$Catalytic reduction$$

$$A = B^1 = B^2$$

$$A = B^2$$

$$A = B^3$$

$$A$$

$$\begin{array}{c|c} NH_2 & NR^1R^2 \\ \hline Het & B^1 = B^2 \text{ ym} \\ \hline \\ NH_2 & NR^1R^2 \\ \hline \\ NH_2 & NR^1R^2 \\ \hline \\ NH_3 & NR^1R^2 \\ \hline \\ NH_4 & NR^1R^2$$

15

wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and nare as defined above.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) can be produced by reacting a heterocyclic dicarboxylic acid imide 5 derivative of the general formula (III-2) with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, in the presence of an inert solvent to obtain a heterocyclic dicarboxylic acid diamide derivative of the general formula (I-3), subjecting said heterocyclic dicarboxylic acid diamide derivative (I-3) to catalytic reduction with hydrogen after or without isolation to obtain a heterocyclic dicarboxylic acid diamide derivative of the general formula (I-2), diazotizing said heterocyclic dicarboxylic acid diamide derivative (I-2) after or without isolation, and then adding a metal salt thereto.

- (4-1). General formula (III-2) \rightarrow general formula (I-3) 20 In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-5).
- (4-2). General formula (I-3) \rightarrow general formula (I-2) In the case of this reaction, the desired 25 compound can be produced in a manner similar to that described in production process (3-3).

(4-3). General formula (I-2) \rightarrow general formula (I-1) In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (3-4).

5 Production process 5.

wherein R^1 , R^2 , R^3 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, m and n are as defined above.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-5) or the general formula (I-4) can be produced by reacting a heterocyclic dicarboxylic acid anhydride of the general formula (VI) with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, in the presence of an inert solvent to 10 obtain a heterocyclic dicarboxylic acid amide of the general formula (IV-3), and treating said heterocyclic dicarboxylic acid amide after or without isolation as follows: when its R2 is a hydrogen atom, the heterocyclic dicarboxylic acid amide (IV-3) is 15 subjected to condensation reaction in the presence of a condensing agent to obtain a compound of the general formula (IX), and said compound (IX) is reacted with an aromatic amine of the general formula (V-1) in the presence of an inert solvent after or without isolating 20 the compound (IX); or when its R2 is other than a hydrogen atom, the heterocyclic dicarboxylic acid amide (IV-3) is condensed with an aromatic amine of the general formula (V-1) in the presence of a condensing agent.

Alternatively, a heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) or the general formula (I-4) can be produced by reacting a heterocyclic dicarboxylic acid anhydride of the general

formula (VI) with an aromatic amine of the general formula (V-1) in the presence of an inert solvent to obtain a heterocyclic dicarboxylic acid amide of the general formula (IV-4), and treating said heterocyclic dicarboxylic acid amide (IV-4) after or without 5 isolation as follows: when its R3 is a hydrogen atom, the heterocyclic dicarboxylic acid amide (IV-4) is subjected to condensation reaction in the presence of a condensing agent to obtain a compound of the general formula (IX-1), and said compound (IX-1) is reacted with an amine or its salt of the general formula (II-1)or the general formula (II-2), respectively, in the presence of an inert solvent after or without isolating the compound (IX-1); or when its R^3 is other than a 15 hydrogen atom, the heterocyclic dicarboxylic acid amide (IV-4) is condensed with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, in the presence of a condensing agent.

(5-1). General formula (VI) \rightarrow general formula (IV-3) 20 or general formula (IV-4)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-3).

General formula (IV-3) \rightarrow general formula (IX) (5-2). or general formula (IV-4) \rightarrow general formula (IX-1) 25 In the case of this reaction, the desired

compound can be produced according to the process described in J. Med. Chem., 10, 982 (1967).

(5-3). General formula (IV-3) or general formula (IV-4) \rightarrow general formula (I-4)

The production can be carried out by reacting a heterocyclic dicarboxylic acid amide derivative of the general formula (IV-3) or the general formula (IV-4) with an amine of the general formula (II-1) or (II-2) or the general formula (V-1) in the presence of a condensing agent and an inert solvent. It is also possible to carry out the reaction in the presence of a base if necessary.

The inert solvent used in the reaction includes, for example, tetrahydrofuran, diethyl ether, dioxane, chloroform and dichloromethane. As the condensing agent used in the reaction, any condensing agent may be used so long as it is used in conventional amide synthesis. The condensing agent includes, for example, Mukaiyama reagent (2-chloro-N-methyl-

pyridinium iodide), DCC (1,3-dicyclohexylcarbodiimide),
CDI (carbonyl diimidazole) and DEPC (diethyl
cyanophosphonate). As to the amount of the condensing
agent used, the condensing agent may be used in an
amount properly chosen in the range of 1 mole to excess
moles per mole of the heterocyclic dicarboxylic acid

amide derivative of the general formula (IV-3) or the general formula (IV-4).

25

The base usable in the reaction includes, for example, organic bases such as triethylamine, pyridine, etc.; and inorganic bases such as potassium carbonate. As to the amount of the base used, the base may be used in an amount properly chosen in the range of 1 mole to excess moles per mole of the heterocyclic dicarboxylic acid amide derivative of the general formula (IV-3) or the general formula (IV-4).

As to the reaction temperature, the reaction

10 may be carried out at 0°C to the boiling point of the

inert solvent used. Although the reaction time is

varied depending on the scale of reaction, the reaction

temperature and the like, it ranges from several

minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystallization, column chromatography, etc., whereby the desired compound can be produced.

(5-4). General formula (IX) \rightarrow general formula (I-5) or general formula (IX-1) \rightarrow general formula (I-1)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-5).

Production process 6.

wherein R, R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and n are as defined above.

A heterocyclic dicarboxylic acid diamide

5 derivative of the general formula (I-1) can be produced
by halogenating a heterocyclic dicarboxylic acid ester
derivative of the general formula (X) in the presence
or absence of an inert solvent to obtain a heterocyclic
dicarboxylic acid halide of the general formula (XI),

10 reacting said heterocyclic dicarboxylic acid halide
(XI) with an aromatic amine of the general formula (V)
in the presence of an inert solvent and a base after or

without isolating the halide (XI), to obtain a
heterocyclic dicarboxylic acid amide of the general
formula (IV-3), hydrolyzing said heterocyclic
dicarboxylic acid amide (IV-3) in the presence or

5 absence of an inert solvent after or without isolating
the amide (IV-3), to obtain a heterocyclic dicarboxylic
acid amide of the general formula (IV-1), subjecting
said heterocyclic dicarboxylic acid amide (IV-1) to
condensation reaction after or without isolation to

10 obtain a heterocyclic dicarboxylic acid isoimide
derivative of the general formula (IX-1), reacting said
heterocyclic dicarboxylic acid isoimide derivative (IX1) with an amine or its salt of the general formula
(II-1) or the general formula (II-2), respectively.

15 (6-1). General formula (X) \rightarrow general formula (XI) As the inert solvent usable in this reaction, any inert solvent may be used so long as it does not markedly inhibit the progress of the reaction. can be exemplified inert solvents including aromatic hydrocarbons such as benzene, toluene, xylene, etc.; 20 halogenated hydrocarbons such as dichloromethane, chloroform, carbon tetrachloride, etc.; chlorinated aromatic hydrocarbons such as chlorobenzene, dichlorobenzene, etc.; acyclic or cyclic ethers such as 25 diethyl ether, dioxane, tetrahydrofuran, etc.; and esters such as ethyl acetate, etc. These inert solvents may be used singly or as a mixture of two or

more thereof.

As a halogenating agent, there can be used halogenating agents such as thionyl chloride, phosphorus oxychloride, phosphorus trichloride, etc.

As to the amount of the halogenating agent used, the halogenating agent may be used in an amount properly chosen in the range of 1 to 10 equivalents per equivalent of the heterocyclic dicarboxylic acid ester derivative of the general formula (X).

As to the reaction temperature, the reaction may be carried out at 0°C to the reflux temperature of the inert solvent used. Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, it may be properly chosen in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystallization, column chromatography, etc., whereby the desired compound can be produced. It is also possible to subject the desired compound to the subsequent reaction without isolation from the reaction system.

(6-2). General formula (XI) → general formula (IV-3)
As the inert solvent usable in this reaction, there can be used, for example, the inert solvents exemplified in production process (1-3).

20

As the base, an inorganic base or an organic base can be used. As the inorganic base, there can be used, for example, hydroxides of alkali metals, such as sodium hydroxide, potassium hydroxide, etc. As the organic base, organic bases such as triethylamine, pyridine and the like can be used. As to the amount of the base used, the base may be used in an amount properly chosen in the range of 0.5 to 3 equivalents per equivalent of the heterocyclic dicarboxylic acid halide of the general formula (XI).

Since the reaction is an equimolar reaction, it is sufficient that the reactants are used in equimolar amounts, though the reaction may be carried out by properly choosing the amount of the aromatic amine of the general formula (V) in the range of 0.5 to 2 equivalents per equivalent of the aromatic dicarboxylic acid halide of the general formula (XI).

As to the reaction temperature, the reaction may be carried out at 0° C to the reflux temperature of the inert solvent used. Although the reaction time is varied depending on the scale of reaction, the reaction temperature and the like, it may be properly chosen in the range of several minutes to 48 hours.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystallization, column chromatography, etc., whereby the

desired compound can be produced. It is also possible to subject the desired compound to the subsequent reaction without isolation from the reaction system.

- (6-3). General formula (IV-3) \rightarrow general formula (IV-1)

 In the case of this reaction, the desired compound can be produced according to production process (1-1).
- (6-5). General formula (IX-1) \rightarrow general formula (I-1)

 In the case of this reaction, the desired compound can be produced according to production

 15 process (I-5).

Production process 7.

Xn
$$CO_2H$$
 $Halogenation$ Het $COhal$ H_2N B^4-B^3 (VII) $(VIII)$ $(VIII$

wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and n are as defined above.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1) can be produced by halogenating a heterocyclic dicarboxylic acid of the general formula (VII) in the presence of a halogenating agent to obtain an acid halide of the general formula (XII), reacting said acid halide (XII) with a

substituted aromatic amine of the general formula (V) in the presence or absence of an inert solvent to obtain a heterocyclic dicarboxylic acid isoimide and a heterocyclic dicarboxylic acid imide of the general

- formulas (IX-1) and (III), respectively, and reacting these compounds with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, in the presence or absence of an inert solvent after or without isolating said compounds.
- 10 (7-1). General formula (VII) \rightarrow general formula (XII)

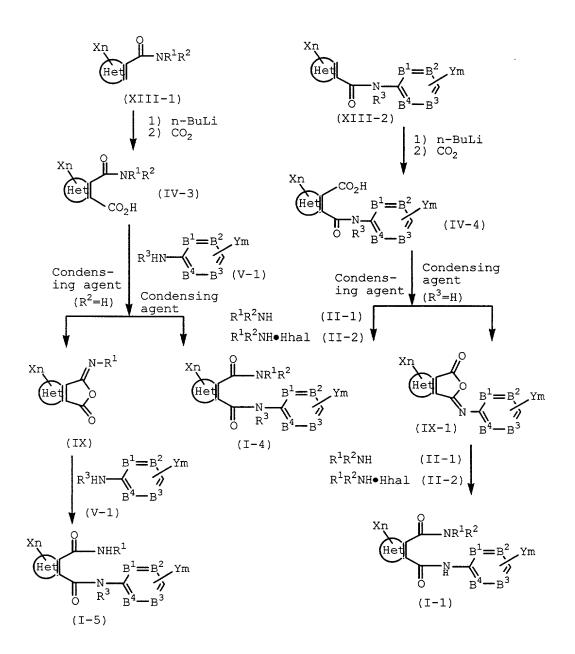
 In the case of this reaction, the desired compound can be produced according to production process (6-1).
- (7-2). General formula (XII) \rightarrow general formulas (IX-1) and (III)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (6-2).

(7-3). General formulas (IX-1) and (III) \rightarrow general 20 formula (I-1)

In the case of this reaction, the desired compound can be produced in a manner similar to that described in production process (1-5).

Production process 8.



wherein R^1 , R^2 , Het, B^1 , B^2 , B^3 , B^4 , X, Y, hal, m and n are as defined above.

A heterocyclic dicarboxylic acid diamide derivative of the general formula (I-1), (I-4) or (I-5)

can be produced by subjecting a heterocyclic carboxylic acid amide of the general formula (XIII-1) or the general formula (XIII-2) to ortho-metallation by the use of a metal reagent such as butyllithium, reacting the ortho-metallation product with carbon dioxide to obtain a heterocyclic dicarboxylic acid amide derivative of the general formula (IV-3) or the general formula (IV-4), and treating this derivative in the same manner as in production processes (5-2) to (5-4).

10 (8-1). General formula (XIII-1) or general formula (XIII-2) \rightarrow general formula (IV-3) or general formula (IV-4)

In the case of this reaction, the production can be carried out by lithiation at the ortho-position according to the method described in J. Org. Chem., 29, 853 (1964) and then introducing carbon dioxide to replace the lithium therewith at -80° C to room temperature.

After completion of the reaction, the desired compound is isolated from the reaction system containing the desired compound by a conventional method, and if necessary, purified by recrystallization, column chromatography, etc., whereby the desired compound can be produced.

Production process 9.

wherein R^1 , R^2 , Het, X, Y, hal, m and n are as defined above.

The reaction is oxidation of the nitrogen atom in the heterocyclic ring. A heterocyclic

5 dicarboxylic acid diamide derivative of the general formula (I-6) can be produced by reacting a heterocyclic dicarboxylic acid imide of the general formula (III-4) with an oxidizing agent in the presence of an inert solvent to obtain a heterocyclic

10 dicarboxylic acid imide derivative of the general formula (III-5), and reacting said heterocyclic dicarboxylic acid imide derivative (III-5) with an amine or its salt of the general formula (II-1) or the general formula (II-2), respectively, after or without

15 isolating the derivative (III-5).

(9-1). General formula (III-4) \rightarrow general formula (III-5)

As the inert solvent usable in this reaction, any inert solvent may be used so long as it does not inhibit the progress of the reaction. There can be exemplified dichloromethane, chloroform, carbon tetrachloride, chlorobenzene, water, acetic acid, ethyl acetate and trifluoroacetic acid. These inert solvents may be used singly or as a mixture of two or more thereof.

As the oxidizing agent used in the reaction, hydrogen peroxide, m-chloroperbenzoic acid and peracetic acid can be exemplified.

As to the reaction temperature for said

15 reaction, the reaction may be carried out at 0°C to

100°C. Although the reaction time is varied depending
on the scale of reaction, the reaction temperature and
the like, it may be properly chosen in the range of
several minutes to 48 hours.

20 (9-2). General formula (III-5) \rightarrow general formula (I-6)

In the case of this reaction, the production

can be carried out according to production process (I-5).

Typical examples of the heterocyclic

25 dicarboxylic acid diamide derivative of the general formula (I) of the present invention are given in

Tables 1 to 33 but they are not intended in any way to

limit the scope of the present invention. General formula (I)

$$\begin{array}{c|c} Xn & Z^1 \\ NR^1R^2 \\ NR^1R^2 \\ NR^3 & B^4 - B^3 \end{array}$$
 (I)

Table 1
$$(Z^1 = Z^2 = 0, R^3 = H, Het = Q1, B^1 = B^2 = B^3 = B^4 = C)$$

No	R¹	R²	р	Xn	Ym	Physical property m.p. °C
1	Н	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
2	СН₃	Н	0	Н	4-CF ₃	
3	СН₃	Н	0	Н	2-CH ₃ -4-Cl	
4	СНз	Н	0	Н	2-CH ₃ -4-OCHF ₂	
5	СНз	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
6	СН₃	Н	0	Н	2-CH ₃ -4-OCF ₃	
7	C ₂ H ₅	Н	0	Н	2-CH ₃ -4-CF ₃	
8	C ₂ H ₅	Н	0	Н	2-CH ₃ -4-OCHF ₂	
9	C ₂ H ₅	Н	0	Н	2-CH ₃ -4-OCF ₃	
10	C ₂ H ₅	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
11	C ₂ H ₅	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 1 (cont'd)

No	R¹	R²	р	Xn	Ym	Physical property
						m. p. °C
12	C ₂ H ₅	C ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	108-111
13	C ₂ H ₅	C ₂ H ₅	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	111-114
14	n-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
15	n-C ₃ H ₇	Н	0	Н	4-CF ₃	
16	n-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₃	
17	n-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
18	C (CH ₃) ₂ CH ₂	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	64-65
	-SCH₃	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
19	n-C ₃ H ₇	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
20	i−C₃H ₇	Н	0	Н	4-CF ₃	
21	i-C ₃ H ₇	Н	0	Н	2-NO ₂	
22	i-C ₃ H ₇	Н	0	Н	4-NO ₂	
23	i-C ₃ H ₇	Н	0	Н	4-F	
24	i-C ₃ H ₇	Н	0	Н	2-CH ₃	
25	i-C ₃ H ₇	Н	0	Н	4-CF ₃	
26	i-C ₃ H ₇	Н	0	Н	3-CF ₃	
27	i−C₃H₁	Н	0	Н	4-CF ₂ CF ₂ CF ₃	
28	i-C ₃ H ₇	Н	0	Н	4-(CF ₂) ₃ CF ₃	
29	i-C ₃ H ₇	Н	0	Н	4-0CF ₃	
30	i-C ₃ H ₇	Н	0	Н	4-OCF ₂ CHFOC ₃ F ₇ -n	
31	i−C₃H ₇	Н	0	Н	3-SCF ₃	

Table 1 (cont'd)

			Ţ	Ι		T
						Physical
No	R ¹	R ²	р	Xn	Ym	property
						m. p. ℃
32	СН (СН₃	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	paste
	-CH ₂ SCH ₃					
33	i−C₃H ₇	Н	0	Н	4-SCH ₂ CF ₃	
34	i-C ₃ H ₇	Н	0	Н	4-SCF ₂ CHF ₂	
35	i-C ₃ H ₇	Н	0	Н	4-S(CF ₂) ₃ CF ₃	
36	i−C₃H ₇	Н	0	Н	4-SCF (CF ₃) ₂	
37	i-C₃H₁	Н	0	Н	4-SCF ₂ CBrF ₂	
38	i-C₃H₁	Н	0	Н	4-SOCF ₂ CBrF ₂	
39	i−C₃H ₇	Н	0	Н	4-S0(CF ₂) ₃ CF ₃	
40	i−C₃H ₇	Н	0	Н	4-S0 ₂ CH ₂ CF ₃	
41	i−C₃H ₇	Н	0	Н	2, 3-Cl ₂	
42	i-C₃H ₇	Н	0	Н	2, 4-Cl ₂	
43	i-C₃H7	Н	0	Н	3, 4-F ₂	
44	i−C₃H ₇	Н	0	н	2, 4-(CH ₃) ₂	
45	i-C₃H ₇	Н	0	Н	2-C1-4-CF ₃	
46	i−C₃H ₇	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
47	i−C₃H ₇	Н	0	Н	2-C1-4-0CF ₃	
48	i−C₃H ₇	Н	0	Н	2-Br-4-0CF ₃	
49	i-C₃H ₇	Н	0	Н	2-CH ₃ -3-C1	
50	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-Cl	
51	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -5-C1	
52	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-Br	

Table 1 (cont'd)

No	R¹	R²	р	Xn	Ym	Physical property m.p. °C
53	i−C₃H ₇	Н	0	Н	2-CH₃-5-F	
54	i−C₃H₁	Н	0	Н	2-CH ₃ -4-CF ₃	167-169
55	i−C₃H₁	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	188-189
56	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
57	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	134-136
58	i-C ₃ H ₇	Н	1	Н	2-CH ₃ -4-CF (CF ₃) ₂	
59	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₃	
60	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
61	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -3-OCF ₂ CHC1F	
62	i-C₃H₁	Н	0	Н	2-CH ₃ -4-OCF ₂ CHC1F	
63	i−C₃H₁	Н	0	Н	2-CH ₃ -4-CF ₂ CBrF ₂	
64	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF ₂ CCl ₂ F	
65	i-C₃H₁	Н	0	Н	2-CH ₃ -4-OCF ₂ CHFCF ₃	
66	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
67	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
68	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-SC ₃ H ₇ -i	
69	i-C₃H ₇	Н	0	Н	2-CH ₃ -4-OCH ₂ OCH ₃	
70	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCH ₂ SCH ₃	
71	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-COOCH ₃	
72	i−C₃H₁	Н	0	Н	2-CH ₃ -4-OCH ₂ COOCH ₃	
73	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-(F ₅ -Ph0)	
74	i−C₃H₁	Н	0	Н	2-CH ₃ -4-(3-CF ₃ -Ph0)	

Table 1 (cont'd)

				· · · · · · · · · · · · · · · · · · ·		
No	R¹	R ²	p	Xn	Ym	Physical property m.p. °C
75	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-(2-C1-4-CF ₃ -Ph0)	ш. р. С
76	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-(4-C1-Ph-CH ₂ 0)	
77	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-(4-C1-PhS)	
78	i-C ₃ H ₇	Н	0	Н	$2-CH_3-4-(5-CF_3-2-Pyr-0)$	
79	i-C ₃ H ₇	H	0	Н	2-CH ₃ -4-(3-C1-	
			0		5-CF ₃ -2-Pyr-0)	
80	i-C ₃ H ₇	H	0	H	$4-(3-C1-5-CF_3-2-Pvr-S)$	
81	i-C ₃ H ₇	H	0	Н	$2-CH_3-4-P=0(0C_2H_5)_2$	
82	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-0P=S (0CH ₃) ₂	
83	i-C ₃ H ₇	Н	0	Н	2-CF ₃ -4-0CHF ₂	
84	i-C ₃ H ₇	Н	0	Н	3, 5-Cl ₂ -4-0CHF ₂	
85	i-C ₃ H ₇	Н	0	Н	3-N=C (CF ₃)-NH-4	
86	i-C ₃ H ₇	Н	0	Н	3-N=C (CF ₃)-N (CH ₃)-4	į
87	i-C ₃ H ₇	Н	0	4-C1	4-C ₄ H ₉ -n	
88	i-C ₃ H ₇	Н	0	4-C1	4-C ₄ H ₉ -t	
89	i-C₃H ₇	Н	0	4-C1	4-CF(CF ₃) ₂	
90	i-C ₃ H ₇	Н	0	4-C1	4-CF ₂ CF ₂ CF ₃	
91	i-C ₃ H ₇	Н	0	4-C1	4-(CF ₂) ₃ CF ₃	
92	i−C₃H ₇	Н	0	4-C1	4-0CHF ₂	
93	i-C₃H₁	Н	0	4-C1	4-OCF ₂ CHFOC ₃ F ₇ -n	
94	i-C₃H₁	н	0	4-C1	4-SCH ₃	
95	i-C₃H₁	Н	0	4-C1	4-SOCH ₃	

Table 1 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
96	i−C₃H ₇	Н	0	4-C1	4-S0 ₂ CH ₃	
97	i-C ₃ H ₇	Н	0	4-C1	4-SCHF ₂	
98	i-C ₃ H ₇	Н	0	4-C1	3−SCF₃	
99	i−C₃H ₇	Н	0	4-C1	3-S0CF ₃	
100	i-C₃H₁	Н	0	4-C1	4-SCH ₂ CF ₃	
101	i-C ₃ H ₇	Н	0	4-C1	4-SCF ₂ CHF ₂	
102	i-C ₃ H ₇	Н	0	4-C1	4-SCF ₂ CBrF ₂	
103	i−C₃H,	Н	0	4-C1	4-SCF (CF ₃) ₂	
104	i-C₃H ₇	Н	0	4-C1	4-S(CF ₂) ₃ CF ₃	
105	i−C₃H ₇	н	0	4-C1	4-SOCF (CF ₃) ₂	
106	i−C₃H₁	Н	0	4-C1	4-S0 ₂ CH ₂ CF ₃	
107	i−C₃H ₇	Н	0	4-C1	4-S0 ₂ CF ₂ CHF ₂	
108	i−C₃H ₇	Н	0	4-C1	4-COCH ₃	
109	i-C₃H₁	Н	0	4-C1	4-Ph	
110	i-C₃H,	Н	0	4-C1	2, 3-Cl ₂	
111	i−C₃H₁	Н	0	4-C1	2, 4-Cl ₂	
112	i−C₃H ₇	Н	0	4-C1	2, 4-F ₂	
113	i−C₃H ₇	Н	0	4-C1	2-C1-4-F	
114	i−C ₃ H ₇	Н	0	4-C1	2-F-4-C1	
115	i−C₃H ₇	Н	0	4-C1	2, 3, 4-F ₃	
116	i−C₃H ₇	Н	0	4-C1	2, 3-(CH ₃) ₂	
117	i-C ₃ H ₇	Н	0	4-C1	2-CH₃-3-Cl	

Table 1 (cont'd)

		I				
	_					Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
118	i-C₃H₁	Н	0	4-C1	2-CH ₃ -4-C1	
119	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -5-C1	
120	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-Br	
121	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-I	
122	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-OCH ₃	1
123	i−C₃H ₇	Н	0	4-C1	2, 3-(CH ₃) ₂ -4-0CH ₃	
124	i-C ₃ H ₇	Н	0	4-C1	2-C1-4-CF ₃	
125	i-C ₃ H ₇	Н	1	4-C1	2-C1-4-CF (CF ₃) ₂	
126	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-CF ₃	
127	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	162-167
128	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CCl ₃	
129	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
130	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
131	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
132	s-C ₄ H ₉	Н	0	4-C1	2-CH ₃ -4-(CF ₂) ₅ CF ₃	
133	i-C ₄ H ₉	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
134	t-C ₄ H ₉	Н	0	4-C1	2-C1-4-0CF ₃	
135	t−C₄H ₉	Н	0	4-C1	2-Br-4-0CF ₃	
136	t-C ₄ H ₉	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	175-180
137	i−C₃H₁	Н	0	4-Br	2-C1-4-CF ₃	
138	i-C₃H₁	Н	0	4-Br	2-C1-4-CF (CF ₃) ₂	
139	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₃	

Table 1 (cont'd)

No	R¹	R ²	þ	Xn	Ym	Physical property m.p. °C
140	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₂ CF ₃	
141	i-C₃H₁	Н	0	4-Br	2-CH ₃ -4-0CF ₂ CCl ₃	
142	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
143	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF(CF ₃) ₂	
144	i−C₃H₁	Н	0	4-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
145	i−C₃H₁	Н	0	4-Br	2-CH ₃ -4-(CF ₂) ₅ CF ₃	
146	i-C ₃ H ₇	Н	0	4-Br	3-C1-4-0CHF ₂	
147	i−C₃H ₇	Н	0	4-Br	2-C1-4-0CF ₃	
148	i−C₃H₁	Н	0	4-Br	2-Br-4-0CF ₃	
149	i-C ₃ H ₇	Н	0	4-Br	2-Br-4-CF(CF ₃) ₂	
150	i−C₃H,	Н	0	4-I	2-CH ₃ -3-C1	
151	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-C1	
152	i−C₃H,	Н	0	4-I	2-CH ₃ -5-C1	
153	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-Br	
154	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-I	
155	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-OCH ₃	
156	i−C₃H ₇	Н	0	4-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
157	i-C₃H₁	Н	0	4-I	2-C1-4-CF ₃	
158	i−C₃H₁	Н	0	4-I	2-C1-4-CF(CF ₃) ₂	
159	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₃	
160	i-C₃H₁	Н	1	4-I	2-CH ₃ -4-CF ₂ CF ₃	
161	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-0CF ₂ CCl ₃	

Table 1 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
162	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
163	i-C₃H,	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
164	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
165	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₅ CF ₃	
166	i−C₃H ₇	Н	0	4-I	3-C1-4-0CHF ₂	
167	i−C₃H₁	Н	0	4-I	2-C1-4-0CF ₃	
168	i−C₃H₁	Н	0	4-I	2-Br-4-0CF ₃	
169	i-C₃H₁	Н	0	4-I	2-Br-4-CF(CF ₃) ₂	
170	i−C₃H ₇	Н	0	6-C1	2-CH ₃ -3-C1	
171	i−C₃H₁	Н	0	6-C1	2-CH ₃ -4-Cl	
172	i−C₃H,	Н	0	6-C1	2-CH ₃ -5-C1	
173	i−C₃H,	Н	0	6-C1	2-CH ₃ -4-Br	
174	i−C₃H₁	Н	0	6-C1	2-CH ₃ -4-I	
175	i−C₃H₁	Н	0	6-C1	2-CH ₃ -4-OCH ₃	
176	i−C₃H₁	Н	0	6-C1	2, 3-(CH ₃) ₂ -4-0CH ₃	
177	i−C₃H ₇	Н	0	6-C1	2-C1-4-CF ₃	
178	i-C ₃ H ₇	Н	0	6-C1	2-C1-4-CF(CF ₃) ₂	
179	i-C₃H ₇	н	0	6-C1	2-CH ₃ -4-CF ₃	
180	i-C ₃ H ₇	Н	0	6-C1	2-CH ₃ -4-CF ₂ CF ₃	
181	i-C₃H₁	Н	0	6-C1	2-CH ₃ -4-OCF ₂ CCl ₃	
182	i−C₃H ₇	Н	0	6-C1	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
183	i-C ₃ H ₇	Н	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	

Table 1 (cont'd)

No	R¹	R ²	р	Xn	Ym	Physical property
						m.p. ℃
184	i−C₃H ₇	Н	0	6-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
185	i-C ₃ H ₇	Н	0	6-C1	2-CH ₃ -4-(CF ₂) ₅ CF ₃	
186	i−C₃H ₇	Н	0	6-C1	3-C1-4-0CHF ₂	
187	i-C ₃ H ₇	Н	0	6-C1	2-C1-4-0CF ₃	
188	i-C ₃ H ₇	Н	0	6-C1	2-Br-4-0CF ₃	
189	i-C ₃ H ₇	Н	0	6-C1	2-Br-4-CF(CF ₃) ₂	
190	i−C₃H ₇	Н	0	6-I	2-CH₃-3-Cl	
191	i-C₃H₁	Н	0	6-I	2-CH₃-4-Cl	
192	i-C ₃ H ₇	Н	0	6-I	2-CH ₃ -5-Cl	
193	i−C₃H ₇	Н	0	6-I	2-CH ₃ -4-Br	
194	i-C₃H,	Н	0	6-I	2-CH ₃ -4-I	
195	i−C₃H,	Н	0	6-I	2-CH ₃ -4-0CH ₃	
196	i−C₃H ₇	Н	0	6-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
197	i−C₃H ₇	Н	0	6-I	2-C1-4-CF ₃	
198	i-C ₃ H ₇	Н	0	6-I	2-C1-4-CF(CF ₃) ₂	
199	i-C ₃ H ₇	Н	0	6-I	2-CH ₃ -4-CF ₃	
200	i−C₃H₁	Н	0	6-I	2-CH ₃ -4-CF ₂ CF ₃	
201	i−C₃H₁	Н	0	6-I	2-CH ₃ -4-OCF ₂ CCl ₃	
202	i−C₃H₁	Н	0	6-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
203	i−C₃H ₇	Н	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
204	i−C₃H ₇	Н	0	6-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
205	i−C₃H ₇	Н	0	6-I	2-CH ₃ -4-(CF ₂) ₅ CF ₃	

Table 1 (cont'd)

.,	p.l	D 2				Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
206	i-C ₃ H ₇	Н	0	6-I	3-C1-4-0CHF ₂	
207	i−C₃H ₇	Н	0	6-I	2-C1-4-0CF ₃	
208	i−C₃H₁	Н	0	6-I	2-Br-4-0CF ₃	
209	i−C₃H ₇	Н	0	6-I	2-Br-4-CF(CF ₃) ₂	
210	i-C₃H₁	Н	0	6-I	2-CH ₃ -4-CF ₃	
211	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₃	
212	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₂ CCl ₃	
213	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
214	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
215	i-C₃H ₇	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
216	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₅ CF ₃	
217	i−C₃H ₇	Н	0	4-I	3-C1-4-OCHF ₂	
218	i-C₃H ₇	Н	0	4-I	2-C1-4-OCF ₃	
219	i-C₃H ₇	Н	0	4-I	2-Br-4-0CF ₃	

Table 2 $(Z^1 = Z^2 = 0, R^3 = H, Het = Q2, B^1 = B^2 = B^3 = B^4 = C)$

				,		
No	R¹	R ²		V	V	Physical
No	K	K	p	Xn	Ym	property
						m.p. ℃
220	i−C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF ₃	
221	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCHF ₂	
222	i-C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₃	
223	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCBrF ₂	
224	i−C₃H₁	Н	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
225	i−C₃H ₇	Н	0	Н	2-CH ₃ -3-OCF ₂ CHC1F	
226	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CCl ₂ F	
227	i-C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CBrF ₂	
228	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
229	i-C₃H₁	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
230	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	234-236
231	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-SCF ₂ CBrF ₂	
232	i−C₃H₁	Н	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
233	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
234	i-C ₃ H ₇	Н	0	Н	2-C1-4-CF ₃	
235	i-C ₃ H ₇	Н	0	Н	2-C1-4-CF(CF ₂) ₂	
236	i-C ₃ H ₇	H	0	Н	3-C1-4-OCHF ₂	
237	i-C ₃ H ₇	Н	0	Н	3-F-4-0CHF ₂	
238	i-C ₃ H ₇	Н	0	Н	2-C1-4-0CF ₃	
239	i-C₃H₁	Н	1	Н	2-CH ₃ -4-CF ₂ CF ₃	
240	i-C₃H₁	Н	1	Н	2-CH ₃ -4-CF(CF ₃) ₂	*

Table 2 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
110	IX :	K	р	AII	¥ 111	m.p. °C
0.41	. 6 11	7.7		0.01	0 CH 4 CD CD	ш. р. С
241	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
242	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	315(dec.)
243	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
244	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
245	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCHF ₂ -5-C1	
246	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CHF ₂ -5-C1	
247	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-SCHF ₂	
248	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-(F ₅ -Ph0)	
249	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-(5-CF ₃ -2	
					-Pyr-0)	
250	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-(3-C1-	
					5-CF ₃ -2-Pyr-0)	
251	i-C ₃ H ₇	Н	0	2-C1	$2-CH_3-4-P=0(0C_2H_5)_2$	
252	i-C₃H₁	Н	0	2-C1	2-CH ₃ -4-OP=S (OCH ₃) ₂	
253	i-C ₃ H ₇	Н	0	2-C1	2-CF ₃ -4-0CHF ₂	
254	i-C₃H ₇	Н	0	2-C1	3-CF ₃ -4-0CHF ₂	
255	i-C ₃ H ₇	Н	0	2-C1	$3-N=C(CF_3)-0-4$	
256	i−C₃H ₇	Н	0	2-C1	3-N=C (CF ₃) -NH-4	
257	i-C ₃ H ₇	Н	0	2-C1	3-N=C(CF ₃)-N(CH ₃)-4	
258	i-C₃H₁	Н	0	2-C1	2-CH ₃ -4-OCF3	229-231
259	i−C₃H₁	Н	0	2-Br	2-CH ₃ -4-CF ₃	

Table 2 (cont'd)

						Physical
No	R ¹	R ²	р	Xn	Ym	property
			P	1111	1111	m.p. °C
260	i-C ₃ H ₇	Н	0	2-Br	2_CU _4_OCUE	ш. р. С
					2-CH ₃ -4-0CHF ₂	
261	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCF ₃	
262	i−C₃H ₇	H	0	2-Br	2-CH ₃ -4-OCBrF ₂	
263	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCF ₂ CHF ₂	
264	i−C ₃ H ₇	Н	0	2-Br	2-CH ₃ -3-OCF ₂ CHC1F	
265	i-C₃H₁	Н	0	2-Br	2-CH ₃ -4-0CF ₂ CCl ₂ F	
266	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCF ₂ CBrF ₂	
267	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
268	i-C₃H₁	Н	0	2-Br	2-CH ₃ -4-CF ₂ CF ₃	
269	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
270	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -4-SCF ₂ CBrF ₂	
271	i−C₃H,	Н	0	2-Br	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
272	i−C₃H₁	Н	0	2-Br	2-CH3-4-SO2CH2CF2CHF2	
273	i−C₃H ₇	Н	0	2-Br	2-C1-4-CF ₃	
274	i-C₃H₁	Н	0	2-Br	2-C1-4-CF(CF ₂) ₂	
275	i−C₃H₁	Н	0	2-Br	3-C1-4-0CHF ₂	
276	i−C₃H₁	Н	0	2-Br	3-F-4-0CHF ₂	
277	i−C₃H₁	Н	0	2-Br	2-C1-4-OCF ₃	
278	i−C₃H ₇	Н	0	2-Br	2-Br-4-0CF ₃	
279	i-C ₃ H ₇	Н	0	2-Br	3, 5-C1 ₂ -4-0CHF ₂	
280	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
281	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	

Table 2 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
			-	William Co.		m.p. ℃
282	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-0CH ₃	
283	i−C₃H ₇	Н	0	2-I	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
284	i-C₃H₁	Н	0	2-I	2, 4-(CH ₃) ₂ -3-0CHF ₂	
285	i-C ₃ H ₇	Н	0	2-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
286	i−C₃H ₇	Н	0	2-I	2-C1-4-0CF ₃	
287	i-C₃H ₇	Н	0	2-I	2-Br-4-0CF ₃	,
288	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCHF ₂	
289	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCF ₃	
290	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCBrF ₂	
291	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCF ₂ CHFCF ₃	
292	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-OCF ₂ CHC1F	
293	i-C₃H ₇	Н	0	2-I	2-CH ₃ -4-0CF ₂ CHF ₂	
294	i-C₃H,	Н	0	2-I	2-CH ₃ -3-C1-4-OCHF ₂	
295	i−C₃H,	Н	0	2-I	2, 3-(CH ₃) ₂ -4-0CHF ₂	
296	i−C₃H,	Н	0	2-I	2-CH ₃ -4-SCH ₃	
297	i−C₃H ₇	Н	0	2-I	2-CH ₃ -4-(3-CF ₃ -Ph0)	
298	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-(3-C1-	
					5-CF ₃ -2-Pyr-0)	
299	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-(5-CF ₃ -2-Pyr-0)	
300	i-C ₃ H ₇	Н	0	5-C1	-3-0CH ₂ 0-4-	
301	i-C ₃ H ₇	Н	0	5-C1	4-CF ₃	
302	i−C₃H ₇	Н	0	5-C1	4-0CF ₃	

Table 2 (cont'd)

						Physical
No	R¹	R ²	p	Xn	Ym	property
						m.p. ℃
303	i-C ₃ H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂	
304	i-C ₃ H ₇	Н	0	5-C1	2, 4, 6-(CH ₃) ₃	
305	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -3-C1	
306	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-C1	
307	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -5-C1	
308	i−C₃H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-Cl	
309	i−C₃H ₇	Н	0	5-C1	2, 4-(CH ₃) ₂ -3-Cl	
310	i−C₃H ₇	Н	0	5-C1	2-C ₂ H ₅ -4-C1	
311	i−C₃H₁	Н	0	5-C1	2-CH ₃ -4-Br	
312	i−C₃H₁	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-Br	
313	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-I	
314	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-F	
315	i-C ₃ H ₇	Н	0	5-C1	2-C1-4-CF ₃	
316	i-C₃H ₇	Н	0	5-C1	2-CH ₃ -4-CF ₃	
317	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
318	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
319	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
320	i−C₃H ₇	Н	0	5-C1	2, 4-(CH ₃) ₂ -3-OCHF ₂	
321	i−C₃H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-0CH ₃	
322	i−C₃H₁	Н	0	5-C1	2-CH ₃ -4-OCH ₃	
323	i−C₃H₁	Н	0	5-C1	2-C1-4-0CF ₃	
324	i−C₃H ₇	Н	0	5-C1	2-Br-4-0CF ₃	

Table 2 (cont'd)

						Physical
No	R¹	\mathbb{R}^2	p	Xn	Ym	property
			-			m. p. ℃
325	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-0CHF ₂	
326	i−C₃H ₇	H	0	5-C1	2-CH ₃ -4-0CF ₃	
327	i-C₃H ₇	H	0	5-C1	2-CH ₃ -4-OCBrF ₂	
328	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHFCF ₃	
329	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHC1F	
330	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-0CF ₂ CHF ₂	
331	i-C ₃ H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-0CHF ₂	
332	i-C₃H₁	Н	0	5-C1	2-CH ₃ -3-C1-4-OCHF ₂	
333	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-SCH ₃	
334	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-(3-CF ₃ -PhO)	
335	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-(5-CF ₃ -2-Pyr-0)	
336	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-(3-C1-5	
					$-CF_3-2-Pyr-0$	
337	i−C₃H ₇	Н	0	5-C1	-3-0CH ₂ 0-4-	
338	i−C₃H ₇	Н	0	5-C1	2-C1-4-0CHF ₂	
339	i−C₃H ₇	Н	0	5-I	4-C1	
340	i−C₃H ₇	Н	0	5-I	4-Br	
341	i−C ₃ H ₇	Н	0	5-I	4-I	
342	i−C₃H ₇	Н	0	5-I	3-CF ₃	
343	i−C₃H ₇	Н	0	5-I	4-CF ₃	
344	i-C ₃ H ₇	Н	0	5-I	4-CF ₂ CF ₂ CF ₃	
345	i−C ₃ H ₇	Н	0	5-I	4-CF (CF ₃) ₂	

Table 2 (cont'd)

	- 1	D 2				Physical
No	R¹	R ²	р	Xn	Ym	property
-						m.p. ℃
346	i−C ₃ H ₇	Н	0	5-I	4-0CF ₃	
347	i-C ₃ H ₇	Н	0	5-I	4-OCF ₂ CHFOCF ₃	
348	i-C ₃ H ₇	Н	0	5-I	4-SCHF ₂	
349	i-C ₃ H ₇	Н	0	5-I	4-SCH ₂ CF ₃	
350	i-C₃H ₇	Н	0	5-I	4-SCF ₂ CHF ₂	
351	i-C ₃ H ₇	Н	0	5-I	4-SCF ₂ CBrF ₂	
352	i-C ₃ H ₇	Н	0	5-I	4-SCF(CF ₃) ₂	
353	i-C ₃ H ₇	Н	0	5-I	4-S (CF ₂) ₃ CF ₃	
354	i-C ₃ H ₇	Н	0	5-I	3, 4-F ₂	
355	i-C ₃ H ₇	Н	0	5–I	2-CH ₃ -3-C1	
356	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-Cl	
357	i-C₃H ₇	Н	0	5-I	2-CH₃-5-C1	
358	i-C ₃ H ₇	Н	0	5-I	2, 4-(CH ₃) ₂ -3-Cl	
359	i−C ₃ H ₇	Н	0	5-I	2, 3-(CH ₃) ₂ -4-C1	
360	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF ₃	
361	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF ₂ CF ₃	
362	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF ₂ CF ₂ CF ₃	
363	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF (CF ₃) ₂	
364	i-C ₃ H ₇	Н	0	5-I	2-C1-4-(CF ₂) ₃ CF ₃	
365	i-C₃H₁	Н	0	5-I	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
366	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-Br	
367	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-I	

Table 2 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
368	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-F	
369	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF ₃	
370	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -3-CF ₃	
371	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-CF ₃	
372	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
373	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
374	i−C₃H,	Н	0	5–I	2-CH ₃ -4-CF(CF ₃) ₂	
375	i-C₃H,	Н	0	5-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
376	i−C₃H₁	Н	0	5-I	2-CH ₃ -4-OCH ₃	
377	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-0-C ₃ H ₇ -i	
378	i-C₃H₁	Н	0	5–I	2, 3-(CH ₃) ₂ -4-0CH ₃	
379	i-C₃H₁	Н	0	5–I	2-CH ₃ -4-OCH ₂ CF ₃	
380	i−C₃H,	Н	0	5–I	2-CH ₃ -4-OCF ₂ CBrF ₂	
381	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CCl ₂ F	
382	i−C₃H ₇	Н	0	5-Br	3-F-4-0CHF ₂	
383	i−C₃H₁	Н	0	5-Br	3, 5-Cl ₂ -4-0CHF ₂	
384	i−C₃H ₇	Н	0	5-Br	3-0CH ₃ -4-0CHF ₂	
385	i−C₃H,	Н	0	5-Br	3, 4-(OCHF ₂) ₂	
386	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-0CF ₃	
387	i−C₃H₁	Н	0	5-Br	2-CH ₃ -4-OCHF ₂	
388	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
389	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCBrF ₂	

Table 2 (cont'd)

1 1					
No R ¹	R^2	p	Xn	Ym	Physical
	1		A11	1 111	property m.p. ℃
390 i-C ₃ H ₂	Н	0	5-Br	2-CH ₃ -4-0CF ₂ CHF ₂	m. p. C
391 i−C₃H;		0	5-Br	2-CH ₃ -4-OCF ₂ CHF ₂ -5-C1	
392 i-C ₃ H ₂		0	5-Br	2-CH ₃ -4-OCF ₂ CHC1F	-
393 i-C ₃ H ₇	1	0	5-Br	2-CH ₃ -4-OCF ₂ CHFCF ₃	
394 i-C ₃ H ₇		0	5-Br	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
395 i-C ₃ H ₇	İ	0	5-Br	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
396 i-C ₃ H ₇		0	5-Br	2-CH ₃ -4-0CHF ₂ -5-C1	
397 i-C ₃ H ₇		0	5-Br	2-CH ₃ -4-(F ₅ -Ph0)	
398 i-C ₃ H ₇		0	5-Br	2-CH ₃ -4-(3-CF ₃ -Ph0)	
399 i-C ₃ H ₇	H	0	5-Br	2-CH ₃ -4-(5-CF ₃ -2-	
				Pyr-0)	
400 i-C ₃ H ₇	Н	0	2−CH₃	2-CH ₃ -4-OCF ₃	
401 i-C ₃ H ₇	Н	0	2-CH ₃	2-CH ₃ -4-CF ₂ CF ₃	
402 i-C ₃ H ₇	H	0	2-CH ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
403 i-C ₃ H ₇	Н	0	2-СН ₃	2-CH ₃ -4-CF (CF ₃) ₂	Į.
404 i-C ₃ H ₇	Н	0	2-CH ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
405 i-C ₃ H ₇	H	0	2-CH ₃	2-C1-4-CF(CF ₃) ₂	
406 i-C ₃ H ₇	Н	0	2-СН3	2-C1-4-(CF ₂) ₃ CF ₃	
407 i-C ₃ H ₇	Н	0	2−СН₃	$2-C_2H_5-4-CF(CF_3)_2$	
408 i-C ₃ H ₇	Н	0	2-CH₃	2-F-4-CF ₂ CF ₃	ì
409 i-C ₃ H ₇	Н	0	2-CF ₃	2-CH ₃ -4-0CF ₃	
410 i-C ₃ H ₇	Н	0	2-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	
			- 0		

Table 2 (cont'd)

No	R¹	R ²	р	Xn	Ym	Physical property m.p. °C
411	i−C₃H₁	Н	0	2-CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
412	i-C ₃ H ₇	Н	0	2-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	
413	i−C₃H ₇	Н	0	2-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
414	i−C₃H₁	Н	0	2-CF ₃	2-C1-4-CF(CF ₃) ₂	
415	i-C ₃ H ₇	Н	0	2-CF ₃	2-C1-4-(CF ₂) ₃ CF ₃	
416	i-C₃H₁	Н	0	2-CF ₃	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
417	i-C ₃ H ₇	Н	0	2-CF3	2-F-4-CF ₂ CF ₃	
418	i−C₃H₁	Н	0	2-CF ₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
419	i−C₃H ₇	Н	0	2-CF ₃	2-Br-4-CF ₂ CF ₃	
420	i-C ₃ H ₇	Н	0	5-СН3	2-CH ₃ -4-0CF ₃	
421	i-C₃H₁	Н	0	5−СН₃	2-CH ₃ -4-CF ₂ CF ₃	
422	i-C ₃ H ₇	Н	0	5-СН3	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
423	i−C₃H ₇	Н	0	5-СН₃	2-CH ₃ -4-CF(CF ₃) ₂	
424	i−C₃H ₇	Н	0	5-СН₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
425	i−C₃H ₇	Н	0	5-СН3	2-C1-4-CF(CF ₃) ₂	
426	i−C₃H ₇	Н	0	5-СН3	2-C1-4-(CF ₂) ₃ CF ₃	
427	i−C₃H ₇	Н	0	5-СН3	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
428	i−C₃H ₇	Н	0	5-СН₃	2-F-4-CF ₂ CF ₃	
429	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-OCF ₃	
430	i−C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	
431	i−C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
432	i−C₃H ₇	Н	0	5–CF ₃	2-CH ₃ -4-CF (CF ₃) ₂	

Table 2 (cont'd)

No	R¹	R²	р	Xn	Ym	Physical property m.p. °C
433	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
434	i−C₃H ₇	Н	0	5-CF₃	2-C1-4-CF(CF ₃) ₂	
435	i−C₃H₁	Н	0	5-CF ₃	2-C1-4-(CF ₂) ₃ CF ₃	
436	i−C₃H₁	Н	0	5-CF ₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
437	i−C₃H₁	Н	0	5−CF₃	$2-C_2H_5-4-CF(CF_3)_2$	
438	i−C₃H₁	Н	0	5−CF₃	2-F-4-CF ₂ CF ₃	
439	i−C₃H₁	Н	0	5-CF ₃	2-Br-4-CF ₂ CF ₃	
440	i−C₃H ₇	Н	0	5−CF₃	2-Br-4-CF(CF ₃) ₂	

Table 3
$$(R^1 = CH(CH_3)CH_2SCH_3, R^3 = H, Z^1 = Z^2 = O, Heta = Q2, B^1 = B^2 = B^3 = B^4 = C)$$

No	R²	р	Xn	Ym	Physical property m.p. °C
441	Н	0	Н	2-CH ₃ -4-0CF ₃	
442	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
443	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
444	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
445	Н	0	Н	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
446	Н	0	Н	2-C1-4-CF(CF ₃) ₂	

Table 3 (cont'd)

No	R²	р	Xn	Ym	Physical property m.p. °C
447	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	-
448	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
449	Н	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
450	Н	0	Н	2-F-4-CF ₂ CF ₃	
451	Н	0	Н	2-Br-4-CF ₂ CF ₃	
452	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
453	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
454	Н	0	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	
455	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
456	Н	0	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
457	Н	0	2-I	2-CH ₃ -4-CF ₂ CF ₃	
458	Н	0	2-I	2-C1-4-CF(CF ₃) ₂	
459	Н	0	5–I	2-CH ₃ -4-CF(CF ₃) ₂	
460	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	

Table 4 $(R^1 = C(CH_3)_2 CH_2 SCH_3, R^3 = H, Z^1 = Z^2 = O,$ He t = Q2, B¹ = B² = B³ = B⁴ = C)

	,				
N.T.	D2		•		Physical
No	R ²	р	Xn	Ym	property
					m.p. ℃
461	Н	0	Н	2-CH ₃ -4-0CF ₃	
462	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
463	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
464	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	148-149
465	Н	0	Н	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
466	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
467	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	
468	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
469	Н	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
470	Н	0	Н	2-F-4-CF ₂ CF ₃	
471	Н	0	Н	2-Br-4-CF ₂ CF ₃	
472	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
473	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
474	Н	0	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	
475	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
476	Н	0	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
477	Н	0	2-I	2-CH ₃ -4-CF ₂ CF ₃	
478	Н	0	2-I	2-C1-4-CF(CF ₃) ₂	
479	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
480	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	

Table 5 $(R^2 = R^3 = H, Z^1 = Z^2 = O, Het = Q2, p = 0,$ $B^1 = B^2 = B^3 = B^4 = C)$

No	R^{1}	Xn	V	Physical
NO	K	An	Ym	property
401	CH (CH) CH CC H	TT	0 01 4 000	m.p. ℃
481	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-OCF ₃	
482	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₃	
483	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
484	C(CH ₃) ₂ CH ₂ SOCH ₃	Н	2-CH ₃ -4-CF(CF ₃) ₂	180-182
485	CH(CH ₃)CH ₂ SC ₂ H ₅	2-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
486	CH(CH ₃)CH ₂ SC ₂ H ₅	2-I	2-C1-4-CF(CF ₃) ₂	
487	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-OCF ₃	
488	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₃	
489	C (CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
490	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
491	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
492	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-I	2-C1-4-CF(CF ₃) ₂	
493	CH(CH3)CH2NHAc	2-C1	2-CH ₃ -4-OCF ₃	
494	CH(CH3)CH2NHAc	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
495	CH(CH ₃)CH ₂ NHAc	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
496	CH(CH₃)CH₂NHAc	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
497	C(CH ₃) ₂ CH ₂ NHAc	2-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
498	C(CH ₃) ₂ CH ₂ NHAc	5-I	2-C1-4-CF(CF ₃) ₂	
499	CH (CH ₃) C ₂ H ₄ OCH ₃	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	
500	CH (CH ₃) C ₂ H ₄ OCH ₃	2-I	2-CH ₃ -4-CF(CH ₃) ₂	
501	C (CH ₃) ₂ C ₂ H ₄ OCH ₃	5-I	2-CH ₃ -4-CF(CF ₃) ₂	

Table 6 $(Z^1 = Z^2 = 0, R^3 = H, Het = Q3, B^1 = B^2 = B^3 = B^4 = C)$

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
502	i-C₃H ₇	Н	0	Н	2-CH ₃ -4-CF ₃	
503	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCHF ₂	
504	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₃	
505	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCBrF ₂	
506	i-C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
507	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -3-OCF ₂ CHC1F	
508	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CCl ₂ F	
509	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CBrF ₂	
510	i-C₃H₁	Н	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
511	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
512	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	206-208
513	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-SCF ₂ CBrF ₂	
514	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
515	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
516	i−C₃H ₇	Н	0	Н	2-C1-4-CF ₃	
517	i−C₃H ₇	Н	0	Н	2-C1-4-CF (CF ₂) ₂	
518	i-C ₃ H ₇	Н	0	Н	3-C1-4-0CHF ₂	
519	i−C₃H ₇	Н	0	Н	3-F-4-0CHF ₂	
520	i−C₃H₁	Н	0	Н	2-C1-4-OCF ₃	
521	i-C₃H₁	Н	1	Н	2-CH ₃ -4-CF ₂ CF ₃	
522	i−C₃H₁	Н	1	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 6 (cont'd)

		Ī		1		
						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. °C
523	i−C₃H₁	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
524	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	275-277
525	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
526	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
527	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-0CHF ₂ -5-C1	
528	i-C ₃ H ₇	Н	0	2-C1	2-CH ₃ -4-OCF ₂ CHF ₂ -5-C1	
529	i−C₃H₁	Н	0	2-C1	2-CH ₃ -4-SCHF ₂	
530	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-(F ₅ -Ph0)	
531	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-(5-CF ₃ -2	
					-Pyr-0)	
532	i−C₃H,	Н	0	2-C1	2-CH ₃ -4-(3-C1-	-
					5-CF ₃ -2-Pyr-0)	
533	i−C₃H₁	Н	0	2-C1	$2-CH_3-4-P=0(0C_2H_5)_2$	
534	i−C₃H₁	Н	0	2-C1	2-CH ₃ -4-0P=S (OCH ₃) ₂	
535	i−C₃H₁	Н	0	2-C1	2-CF ₃ -4-OCHF ₂	
536	i-C₃H₁	Н	0	2-C1	3-CF ₃ -4-OCHF ₂	
537	i−C₃H ₇	Н	0	2-C1	3-N=C(CF ₃)-0-4	
538	i−C₃H ₇	Н	0	2-C1	3-N=C(CF ₃)-NH-4	
540	i−C₃H ₇	Н	0	2-C1	3-N=C(CF ₃)-N(CH ₃)-4	
541	i−C₃H ₇	Н	0	2-C1	2-CH ₃ -4-OCHF ₂	
542	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-CF ₃	
543	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCHF ₂	

Table 6 (cont'd)

	_	_				Physical
No	R¹	R²	р	Xn	Ym	property
						m. p. ℃
544	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-0CF ₃	
545	i-C₃H₁	Н	0	2-Br	2 -CH $_3$ - 4 -OCBrF $_2$	
546	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -4-OCF ₂ CHF ₂	
547	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -3-OCF ₂ CHC1F	
548	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCF ₂ CCl ₂ F	
549	i-C ₃ H ₇	Н	0	2-Br	2 -CH $_3$ - 4 -OCF $_2$ CBrF $_2$	
550	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
551	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-CF ₂ CF ₃	
552	i−C₃H ₇	Н	0	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
553	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -4-SCF ₂ CBrF ₂	
554	i-C ₃ H ₇	Н	0	2-Br	2-CH3-4-SCH2CF2CHF2	
555	i-C ₃ H ₇	Н	0	2-Br	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
556	i-C ₃ H ₇	Н	0	2-Br	2-C1-4-CF ₃	
557	i-C ₃ H ₇	Н	0	2-Br	2-C1-4-CF(CF ₂) ₂	
558	i-C ₃ H ₇	Н	0	2-Br	3-C1-4-OCHF ₂	
559	i-C ₃ H ₇	Н	0	2-Br	3-F-4-0CHF ₂	
560	i−C₃H₁	Н	0	2-Br	2-C1-4-OCF ₃	
561	i−C₃H ₇	Н	0	2-Br	2-Br-4-0CF ₃	
562	i−C₃H ₇	Н	0	2-Br	3, 5-Cl ₂ -4-OCHF ₂	
563	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
564	i−C₃H,	Н	0	2-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
565	i-C₃H₁	Н	0	2-I	2-CH ₃ -4-0CH ₃	

Table 6 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
			-			m.p. °C
566	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
567	i-C ₃ H ₇	Н	0	2-I	2, 4-(CH ₃) ₂ -3-0CHF ₂	
568	i−C₃H ₇	Н	0	2-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
569	i−C₃H₁	Н	0	2-I	2-C1-4-0CF ₃	
570	i−C₃H₁	Н	0	2-I	2-Br-4-0CF ₃	
571	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-OCHF ₂	
572	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-OCF ₃	
573	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCBrF ₂	
574	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-OCF ₂ CHFCF ₃	
575	i−C₃H₁	Н	0	2-I	2-CH ₃ -4-OCF ₂ CHC1F	
576	i−C₃H ₇	Н	0	2-I	2-CH ₃ -4-OCF ₂ CHF ₂	
577	i−C₃H₁	Н	0	2-I	2-CH ₃ -3-C1-4-0CHF ₂	
578	i−C₃H ₇	Н	0	2-I	2, 3-(CH ₃) ₂ -4-0CHF ₂	
579	i−C₃H ₇	Н	0	2-I	2-CH ₃ -4-SCH ₃	
580	i-C ₃ H ₇	Н	0	2-I	2-CH ₃ -4-(3-CF ₃ -Ph0)	
581	i−C₃H ₇	Н	0	2-I	2-CH ₃ -4-(3-C1-	
					5-CF ₃ -2-Pyr-0)	
582	i−C₃H₁	Н	0	5-C1	2-CH ₃ -4-(5-CF ₃ -2-Pyr-0)	
583	i-C ₃ H ₇	Н	0	5-C1	-3-0CH ₂ 0-4-	
584	i-C₃H₁	Н	0	5-C1	4-CF ₃	
585	i−C₃H₁	Н	0	5-C1	4-0CF ₃	
586	i−C₃H₁	Н	0	5-C1	2, 3-(CH ₃) ₂	
			<u> </u>			

Table 6 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
587	i−C₃H ₇	Н	0	5-C1	2, 4, 6-(CH ₃) ₃	
588	i−C₃H₁	Н	0	5-C1	2-CH ₃ -3-Cl	
589	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-C1	
590	i-C₃H₁	Н	0	5-C1	2-CH₃-5-Cl	
591	i-C ₃ H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-Cl	
592	i−C₃H ₇	Н	0	5-C1	2, 4-(CH ₃) ₂ -3-C1	
593	i-C₃H₁	Н	0	5-C1	2-C ₂ H ₅ -4-Cl	
594	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-Br	
595	i−C₃H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-Br	
596	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-I	
597	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-F	
598	i−C₃H₁	Н	0	5-C1	2-C1-4-CF ₃	
599	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-CF ₃	
600	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
601	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
602	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
603	i-C ₃ H ₇	Н	0	5-C1	2, 4-(CH ₃) ₂ -3-OCHF ₂	
604	i-C ₃ H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-0CH ₃	
605	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-OCH ₃	
606	i-C ₃ H ₇	Н	0	5-C1	2-C1-4-OCF ₃	
607	i-C ₃ H ₇	Н	0	5-C1	2-Br-4-OCF ₃	
608	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-0CHF ₂	

Table 6 (cont'd)

N	n l	\mathbb{D}^2		v	V	Physical
No	R¹	R ²	р	Xn	Ym	property m.p. ℃
609	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-0CF ₃	ш. р. С
610	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCBrF ₂	
611	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHFCF ₃	
612	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHC1F	
613	i-C₃H₁	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHF ₂	
614	i-C₃H ₇	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-0CHF ₂	
615	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -3-C1-4-0CHF ₂	
616	i-C ₃ H ₇	Н	0	5-C1	2−CH₃−4−SCH₃	
617	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-(3-CF ₃ -Ph0)	
618	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-(5-CF ₃ -2-Pyr-0)	
619	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-(3-C1-5	
					$-CF_3-2-Pyr-0)$	
620	i-C₃H₁	Н	0	5-C1	-3-0CH ₂ 0-4-	
621	i-C ₃ H ₇	Н	0	5-C1	2-C1-4-OCHF ₂	
622	i−C₃H ₇	H	0	5-I	4-C1	
623	i-C₃H₁	Н	0	5-I	4-Br	
624	i−C₃H ₇	Н	0	5-I	4-I	
625	i−C₃H ₇	Н	0	5-I	3-CF ₃	
626	i-C ₃ H ₇	Н	0	5-I	4-CF ₃	
627	i−C₃H ₇	Н	0	5-I	4-CF ₂ CF ₂ CF ₃	
628	i-C ₃ H ₇	Н	0	5-I	4-CF(CF ₃) ₂	
629	i-C ₃ H ₇	Н	0	5-I	4-0CF ₃	

Table 6 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
630	i-C ₃ H ₇	Н	0	5-I	4-OCF ₂ CHFOCF ₃	
631	i-C ₃ H ₇	Н	0	5-I	4-SCHF ₂	
632	i-C ₃ H ₇	Н	0	5-I	4-SCH ₂ CF ₃	
633	i-C ₃ H ₇	Н	0	5-I	4-SCF ₂ CHF ₂	
634	i-C ₃ H ₇	Н	0	5-I	4-SCF ₂ CBrF ₂	
635	i-C ₃ H ₇	Н	0	5-I	4-SCF (CF ₃) ₂	
636	i-C ₃ H ₇	Н	0	5-I	4-S (CF ₂) ₃ CF ₃	
637	i−C₃H ₇	Н	0	5-I	3, 4-F ₂	
638	i−C₃H ₇	Н	0	5-I	2-CH ₃ -3-C1	
639	i−C₃H,	Н	0	5-I	2-CH ₃ -4-Cl	
640	i−C₃H ₇	Н	0	5-I	2-CH ₃ -5-C1	
641	i-C ₃ H ₇	Н	0	5-I	2, 4-(CH ₃) ₂ -3-Cl	
642	i-C ₃ H ₇	Н	0	5-I	2, 3-(CH ₃) ₂ -4-Cl	
643	i−C₃H₁	Н	0	5-I	2-C1-4-CF ₃	
644	i-C ₃ H ₇	Н	0	5-I	2-C1-4-CF ₂ CF ₃	
645	i−C₃H ₇	Н	0	5-I	2-C1-4-CF ₂ CF ₂ CF ₃	
646	i−C₃H ₇	Н	0	5-I	2-C1-4-CF(CF ₃) ₂	
647	i−C₃H ₇	Н	0	5-I	2-C1-4-(CF ₂) ₃ CF ₃	
648	i−C₃H ₇	Н	0	5-I	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
649	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-Br	
650	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-I	
651	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-F	

Table 6 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
652	i−C₃H₁	Н	0	5-I	2-C1-4-CF ₃	
653	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -3-CF ₃	
654	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF ₃	
655	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
656	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
657	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
658	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
659	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-OCH ₃	
660	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-0-C ₃ H ₇ -i	
661	i-C ₃ H ₇	Н	0	5-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
662	i-C ₃ H ₇	Н	0	5-I	2-CH ₃ -4-OCH ₂ CF ₃	
663	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-OCF ₂ CBrF ₂	
664	i-C ₃ H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CCl ₂ F	
665	i-C ₃ H ₇	Н	0	5-Br	3-F-4-0CHF ₂	
666	i−C₃H ₇	Н	0	5-Br	3, 5-Cl ₂ -4-OCHF ₂	
667	i-C ₃ H ₇	Н	0	5-Br	3-0CH ₃ -4-0CHF ₂	
668	i-C ₃ H ₇	Н	0	5-Br	3, 4-(0CHF ₂) ₂	
669	i-C ₃ H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₃	
670	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCHF ₂	
671	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
672	i-C₃H₁	Н	0	5-Br	2-CH ₃ -4-OCBrF ₂	
673	i-C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CHF ₂	

Table 6 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
674	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CHF ₂ -5-C1	
675	i-C ₃ H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CHC1F	
676	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CHFCF ₃	
677	i−C₃H₁	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
678	i−C₃H₁	Н	0	5-Br	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
679	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-OCHF ₂ -5-C1	
680	i−C₃H ₇	Н	0	5-Br	$2-CH_3-4-(F_5-Ph0)$	
681	i-C ₃ H ₇	Н	0	5-Br	2-CH ₃ -4-(3-CF ₃ -Ph0)	
682	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-(5-CF ₃ -2-	
					Pyr-0)	
683	i-C ₃ H ₇	Н	0	2-СН3	2-CH ₃ -4-0CF ₃	
684	i-C ₃ H ₇	Н	0	2-СН3	2-CH ₃ -4-CF ₂ CF ₃	
685	i-C ₃ H ₇	Н	0	2-СН3	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
686	i-C ₃ H ₇	Н	0	2−СН₃	2-CH ₃ -4-CF(CF ₃) ₂	
687	i-C₃H ₇	Н	0	2-СН₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
688	i−C₃H ₇	Н	0	2−СН₃	2-C1-4-CF(CF ₃) ₂	
689	i-C ₃ H ₇	Н	0	2-СН3	2-C1-4-(CF ₂) ₃ CF ₃	
690	i-C₃H₁	Н	0	2-СН3	2-C ₂ H ₅ -4-CF (CF ₃) ⁻ ₂	
691	i-C ₃ H ₇	Н	0	2-СН3	2-F-4-CF ₂ CF ₃	
692	i-C ₃ H ₇	Н	0	2-CF ₃	2-CH ₃ -4-OCF ₃	
693	i-C₃H₁	Н	0	2-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	
694	i-C ₃ H ₇	Н	0	2-CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	

Table 6 (cont'd)

No	R¹	R²	р	Xn	Ym	Pyysical property
695	i-C ₃ H ₇	Н	0	9_CE	9 CH 4 CE (CE)	m.p. °C
				2-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	
696	i-C ₃ H ₇	H	0	2-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
697	i−C₃H ₇	Н	0	2-CF ₃	2-C1-4-CF(CF ₃) ₂	
698	i−C₃H₁	Н	0	2-CF ₃	2-C1-4-(CF ₂) ₃ CF ₃	
699	i−C₃H ₇	Н	0	2-CF ₃	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
700	i-C₃H₁	Н	0	2-CF ₃	2-F-4-CF ₂ CF ₃	
701	i−C₃H ₇	Н	0	2-CF ₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
702	i−C₃H ₇	Н	0	2-CF ₃	2-Br-4-CF ₂ CF ₃	
703	i−C₃H₁	Н	0	5−СН₃	2-CH ₃ -4-OCF ₃	
704	i-C ₃ H ₇	Н	0	5-СН3	2-CH ₃ -4-CF ₂ CF ₃	
705	i−C₃H ₇	Н	0	5-СН3	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
706	i−C₃H ₇	Н	0	5-СН3	2-CH ₃ -4-CF(CF ₃) ₂	
707	i−C₃H ₇	Н	0	5-СН₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
708	i−C₃H ₇	Н	0	5−СН₃	2-C1-4-CF(CF ₃) ₂	
709	i−C₃H₁	Н	0	5-СН3	2-C1-4-(CF ₂) ₃ CF ₃	
710	i−C₃H ₇	Н	0	5-СН3	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
711	i−C₃H ₇	Н	0	5-СН3	2-F-4-CF ₂ CF ₃	
712	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-OCF ₃	
713	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	
714	i−C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
715	i−C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	
716	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	

Table 6 (cont'd)

No	R¹	R²	р	Xn	Ym	Physical property m.p. °C
717	i-C ₃ H ₇	Н	0	5−CF₃	2-C1-4-CF(CF ₃) ₂	
718	i-C₃H₁	Н	0	5-CF₃	2-C1-4-(CF ₂) ₃ CF ₃	
719	i−C₃H₁	Н	0	5-CF ₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
720	i-C ₃ H ₇	Н	0	5-CF ₃	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
721	i−C₃H ₇	Н	0	5-CF ₃	2-F-4-CF ₂ CF ₃	
722	i-C₃H₁	Н	0	5-CF ₃	2-Br-4-CF ₂ CF ₃	
723	i−C₃H ₇	Н	0	5-CF ₃	2-Br-4-CF(CF ₃) ₂	

Table 7
$$(R^1 = CH(CH_3)CH_2SCH_3, R^3 = H, Z^1 = Z^2 = O, Heta = Q3, B^1 = B^2 = B^3 = B^4 = C)$$

No	R²	р	Xn	Ym	Physical property m.p. °C
724	Н	0	Н	2-CH ₃ -4-0CF ₃	
725	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
726	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
727	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
728	Н	0	Н	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
729	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
730	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	

Table 7 (cont'd)

					
					Physical
No	R ²	p	Xn	Ym	property
					m.p. ℃
731	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
732	Н	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
733	Н	0	Н	2-F-4-CF ₂ CF ₃	
734	Н	0	Н	2-Br-4-CF ₂ CF ₃	
735	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
736	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
737	Н	0	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	237-239
738	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
739	Н	0	2-1	2-CH ₃ -4-CF(CF ₃) ₂	
740	Н	0	2-1	2-CH ₃ -4-CF ₂ CF ₃	
741	Н	0	2-I	2-C1-4-CF(CF ₃) ₂	
742	Н	0	5-I	2-CH ₃ -4-CF (CF ₃) ₂	
743	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
		1	<u> </u>	J	

Table 8 $(R^1 = C(CH_3)_2 CH_2 SCH_3, R^3 = H, Z^1 = Z^2 = O, Het = Q3, B^1 = B^2 = B^3 = B^4 = C)$

					Physical
No	R ²	р	Xn	Ym	property
					m.p. ℃
744	Н	0	Н	2-CH ₃ -4-OCF ₃	
745	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
746	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
747	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
748	Н	0	Н	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
749	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
750	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	
751	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
752	Н	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
753	Н	0	Н	2-F-4-CF ₂ CF ₃	
754	Н	0	Н	2-Br-4-CF ₂ CF ₃	
755	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
756	Н	0	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
757	Н	0	2-C1	2-CH ₃ -4-CF (CF ₃) ₂	
758	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
759	Н	0	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
760	Н	0	2-I	2-CH ₃ -4-CF ₂ CF ₃	
761	Н	0	2-I	2-C1-4-CF(CF ₃) ₂	
762	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
763	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	

Table 9 $(R^2 = R^3 = H, Z^1 = Z^2 = O, Het = Q3, p = 0,$ $B^1 = B^2 = B^3 = B^4 = C)$

No	R¹	Xn	Ym	Physical property m.p. °C
764	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-0CF ₃	
765	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₃	
766	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
767	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF(CF ₃) ₂	
768	CH(CH ₃)CH ₂ SC ₂ H ₅	2-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
769	CH(CH ₃)CH ₂ SC ₂ H ₅	2-I	2-C1-4-CF(CF ₃) ₂	
770	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-OCF ₃	
771	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₃	
772	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
773	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
774	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
775	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	5-I	2-C1-4-CF(CF ₃) ₂	
776	CH(CH3)CH2NHAc	2-C1	2-CH ₃ -4-OCF ₃	
777	CH(CH₃)CH₂NHAc	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
778	CH(CH₃)CH₂NHAc	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
779	CH(CH₃)CH₂NHAc	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
780	C(CH ₃) ₂ CH ₂ NHAc	2-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
781	C(CH ₃) ₂ CH ₂ NHAc	5-I	2-C1-4-CF(CF ₃) ₂	
782	CH(CH₃)C₂H₄OCH₃	2-C1	2-CH ₃ -4-CF(CF ₃) ₂	
783	CH(CH ₃)C ₂ H ₄ OCH ₃	2-I	2-CH ₃ -4-CF(CH ₃) ₂	
784	C (CH ₃) ₂ C ₂ H ₄ OCH ₃	5-I	2-CH ₃ -4-CF(CF ₃) ₂	

Table 10 $(Z^1 = Z^2 = 0, R^3 = H, Het = Q4, B^1 = B^2 = B^3 = B^4 = C)$

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
785	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF ₃	184-185
786	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCHF ₂	
787	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₃	
788	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCBrF ₂	
789	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
790	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -3-OCF ₂ CHC1F	
791	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CCl ₂ F	
792	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-OCF ₂ CBrF ₂	
793	i-C₃H₁	Н	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
794	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	paste
795	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	159-161
796	i−C₃H ₇	Н	0	Н	2-CH ₃ -4-SCF ₂ CBrF ₂	
797	i-C₃H₁	Н	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
798	i-C₃H₁	Н	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
799	i−C₃H ₇	Н	0	Н	2-C1-4-CF ₃	
800	i-C₃H ₇	Н	0	Н	2-C1-4-CF(CF ₂) ₂	
801	i-C₃H₁	Н	0	Н	3-C1-4-0CHF ₂	
802	i-C₃H,	Н	0	Н	3-F-4-0CHF ₂	
803	i-C₃H₁	Н	1	Н	2-CH ₃ -4-CF ₂ CF ₃	
804	i-C₃H₁	Н	1	Н	2-CH ₃ -4-CF(CF ₃) ₂	108-110
805	C ₂ H ₅	C_2H_5	0	Н	2-CH ₃ -4-CF ₂ CF ₃	130-132

Table 10 (cont'd)

						Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
806	n-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	
807	с-С _з Н ₅	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
808	n-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CBrFCF ₃	
809	i-C₄H ₉	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CHFOCF ₃	
810	i-C ₄ H ₉	Н	0	4-C1	2-CH ₃ -4-OCHF ₂ -5-C1	
811	n-C ₄ H ₉	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CHF ₂ -5-C1	
812	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-SCHF ₂	
813	i-C₃H₁	Н	0	4-C1	2-CH ₃ -4-(F ₅ -Ph0)	
814	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-(5-CF ₃ -2	
					-Pyr-0)	
815	i-C₃H ₇	Н	0	4-C1	2-CH ₃ -4-(3-C1-	
					5-CF ₃ -2-Pyr-0)	
816	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-P=0(0C ₂ H ₅) ₂	
817	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-OP=S (OCH ₃) ₂	
818	i−C₃H ₇	Н	0	4-C1	2-CF ₃ -4-0CHF ₂	
819	i−C₃H ₇	Н	0	4-C1	3-CF ₃ -4-0CHF ₂	
820	i-C₃H₁	Н	0	4-C1	3-N=C(CF ₃)-0-4	
821	i-C₃H₁	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	149-152
822	i-C₃H₁	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
823	с-СзН5	Н	0	4-C1	2-CH ₃ -4-OCHF ₂	
824	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-CF ₃	
825	i-C₃H₁	Н	0	4-C1	2-CH ₃ -4-OCHF ₂	

Table 10 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
826	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-OCF ₃	
827	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-OCBrF ₂	
828	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -4-0CF ₂ CHF ₂	
829	i-C ₃ H ₇	Н	0	4-C1	2-CH ₃ -3-OCF ₂ CHC1F	
830	i−C₃H ₇	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CCl ₂ F	
831	i-C₃H,	Н	0	4-C1	2-CH ₃ -4-OCF ₂ CBrF ₂	
832	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
833	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₂ CF ₃	
834	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-CF(CF ₃) ₂	
835	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-SCF ₂ CBrF ₂	
836	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
837	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
838	i−C₃H ₇	Н	0	4-Br	2-C1-4-CF ₃	
839	i−C₃H ₇	Н	0	4-Br	2-C1-4-CF(CF ₂) ₂	
840	i−C₃H ₇	Н	0	4-Br	3-C1-4-OCHF ₂	
841	i-C ₃ H ₇	Н	0	4-Br	3-F-4-0CHF ₂	
842	i−C₃H ₇	Н	0	4-Br	2-C1-4-OCF ₃	
843	i−C₃H ₇	Н	0	4-Br	2-Br-4-0CF ₃	
844	i-C₃H₁	Н	0	4-Br	3, 5-C1 ₂ -4-0CHF ₂	
845	i−C₃H₁	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
846	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
847	i-C₃H ₇	Н	0	4-I	2-CH ₃ -4-OCH ₃	

Table 10 (cont'd)

						Physical
No	R¹	R ²	p	Xn	Ym	property
						m.p. °C
848	i−C₃H ₇	Н	0	4-I	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
849	i-C ₃ H ₇	Н	0	4-I	2, 4-(CH ₃) ₂ -3-0CHF ₂	
850	i-C ₃ H ₇	Н	0	4-I	2, 3-(CH ₃) ₂ -4-0CH ₃	
851	i-C ₃ H ₇	Н	0	4-I	2-C1-4-0CF ₃	
852	i-C ₃ H ₇	Н	0	4-I	2-Br-4-0CF ₃	
853	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCHF ₂	
854	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₃	
855	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCBrF ₂	
856	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₂ CHFCF ₃	
857	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₂ CHC1F	
858	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₂ CHF ₂	
859	i-C₃H₁	Н	0	4-I	2-CH ₃ -3-C1-4-0CHF ₂	
860	i−C₃H₁	Н	0	4-I	2, 3-(CH ₃) ₂ -4-OCHF ₂	
861	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-SCH ₃	
862	i-C₃H ₇	Н	0	4-I	2-CH ₃ -4-(3-CF ₃ -Ph0)	
863	i−C₃H,	Н	0	4-I	2-CH ₃ -4-(3-C1-	
					5-CF ₃ -2-Pyr-0)	
864	i−C₃H₁	Н	0	5-C1	2-C1-4-CF ₃	
865	i−C₃H₁	Н	0	5-C1	2-CH ₃ -4-CF ₃	
866	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
867	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
868	i-C ₃ H ₇	Н	0	5-C1	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	

Table 10 (cont'd)

	_ 1					Physical
No	R ¹	R²	р	Xn	Ym	property
						m.p. °C
869	i-C₃H₁	Н	0	5-C1	2, 4-(CH ₃) ₂ -3-0CHF ₂	
870	i−C₃H₁	Н	0	5-C1	2, 3-(CH ₃) ₂ -4-0CH ₃	
871	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCH ₃	
872	i-C ₃ H ₇	Н	0	5-C1	2-C1-4-0CF ₃	
873	i−C₃H ₇	Н	0	5-C1	2-Br-4-0CF ₃	
874	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCHF ₂	
875	i−C₃H ₇	Н	0	5-C1	2-CH ₃ -4-OCF ₃	
876	i−C₃H ₇	Н	0	5-C1	2-CH3-4-OCBrF2	
877	i-C₃H,	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHFCF ₃	
878	i-C ₃ H ₇	Н	0	5-C1	2-CH₃-4-OCF₂CHC1F	
879	i−C₃H₁	Н	0	5-C1	2-CH ₃ -4-OCF ₂ CHF ₂	
880	i−C₃H₁	Н	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
881	i−C₃H₁	Н	0	6-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
882	i−C₃H ₇	Н	1	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
883	i-C₃H₁	Н	1	6-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
884	i−C₃H ₇	Н	0	6-C1	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
885	i−C₃H₁	Н	0	4-I	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
886	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-CF (CF ₃) ₂ F ₂	
887	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
888	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
889	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-(CF ₂) ₃ CF	
890	i−C₃H ₇	Н	0	4-I	2-Br-4-0CF ₃	

Table 10 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. °C
891	i-C₃H₁	Н	0	4-I	2-CH ₃ -4-OCHF ₂	
892	i-C₃H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₃	
893	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCBrF ₂	
894	i-C ₃ H ₇	Н	0	4-I	2-CH3-4-OCF2CHFCF3	
895	i-C ₃ H ₇	Н	0	4-I	2-CH ₃ -4-OCF ₂ CHC1F	
896	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
897	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
898	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-CF (CF ₃) ₂	
899	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-SCH ₃	
900	i-C₃H₁	Н	0	5-I	2-CH ₃ -4-(3-CF ₃ -Ph0)	
901	i−C₃H₁	Н	0	5-I	2-CH ₃ -4-(5-CF ₃ -2-Pyr-0)	
902	i−C₃H ₇	Н	0	5-I	2-CH ₃ -4-(3-C1-5	
					-CF ₃ -2-Pyr-0)	,
903	i−C₃H₁	Н	0	6-I	2-CH ₃ -4-CF (CF ₃) ₂	
904	i−C₃H₁	Н	0	6-I	2-C1-4-CF ₂ CF ₃	
905	i−C₃H₁	Н	0	4-Br	2-C1-4-CF ₃	
906	i−C₃H₁	Н	0	4-Br	2-C1-4-CF ₂ CF ₃	
907	i−C₃H₁	Н	0	4-Br	2-C1-4-CF ₂ CF ₂ CF ₃	
908	i−C₃H₁	Н	0	4-Br	2-C1-4-CF(CF ₃) ₂	ļ
909	i−C₃H ₇	Н	0	4-Br	2-C1-4-(CF ₂) ₃ CF ₃	
910	i−C₃H ₇	Н	0	4-Br	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
911	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-Br	

Table 10 (cont'd)

						Physical
No	R¹	R²	5	Xn	Ym	-
NO	K	K	р	All	1111	property
						m.p. °C
912	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-I	
913	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-F	
914	i-C ₃ H ₇	Н	0	4-Br	2-C1-4-CF ₃	
916	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₃	
917	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₂ CF ₃	
918	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
919	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-CF(CF ₃) ₂	
920	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
921	i−C₃H ₇	Н	0	4-Br	2-CH ₃ -4-OCH ₃	
922	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-0-C ₃ H ₇ -i	
923	i-C ₃ H ₇	Н	0	4-Br	2, 3-(CH ₃) ₂ -4-0CH ₃	
924	i−C₃H₁	Н	0	4-Br	2-CH ₃ -4-OCH ₂ CF ₃	
925	i-C ₃ H ₇	Н	0	4-Br	2-CH ₃ -4-OCF ₂ CBrF ₂	
926	i−C₃H₁	Н	0	5-Br	2-CH ₃ -4-OCF ₃	
927	i−C₃H₁	Н	0	5-Br	2-CH ₃ -4-CF ₂ CF ₃	
928	i-C₃H₁	Н	0	5-Br	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
929	i−C₃H ₇	Н	0	5-Br	2-CH ₃ -4-CF(CF ₃) ₂	
930	i-C ₃ H ₇	Н	0	5-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
931	i-C₃H₁	Н	0	5-Br	2-C1-4-CF (CF ₃) ₂	
932	i-C ₃ H ₇	Н	0	5-Br	2-C1-4-(CF ₂) ₃ CF ₃	
933	i-C ₃ H ₇	Н	0	5-Br	2-C ₂ H ₅ -4-CF(CF ₃) ₂	

Table 10 (cont'd)

						Physical
No	R¹	R ²	р	Xn	Ym	property
						m.p. ℃
934	i-C ₃ H ₇	Н	0	5-Br	2-F-4-CF ₂ CF ₃	
935	i-C ₃ H ₇	Н	0	6-Br	2-CH ₃ -4-OCF ₃	
936	i-C ₃ H ₇	Н	0	6-Br	2-CH ₃ -4-CF ₂ CF ₃	
937	i-C ₃ H ₇	Н	0	6-Br	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
938	i-C ₃ H ₇	Н	0	6-Br	2-CH ₃ -4-CF(CF ₃) ₂	
939	i-C ₃ H ₇	Н	0	6-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
940	i-C ₃ H ₇	Н	0	6-Br	2-C1-4-CF(CF ₃) ₂	
941	i-C ₃ H ₇	Н	0	6-Br	2-C1-4-(CF ₂) ₃ CF ₃	
942	i-C ₃ H ₇	Н	0	6-Br	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
943	i-C ₃ H ₇	Н	0	6-Br	2-F-4-CF ₂ CF ₃	
944	i-C₃H ₇	Н	0	6-Br	2-C ₂ H ₅ -4-CF ₂ CF ₃	
945	i−C₃H ₇	Н	0	4-СН3	2-CH ₃ -4-OCF ₃	
946	i-C₃H ₇	Н	0	4−CH ₃	2-CH ₃ -4-CF ₂ CF ₃	
947	i-C₃H ₇	Н	0	4-CH ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
948	i-C ₃ H ₇	Н	0	4-СН3	2-CH ₃ -4-CF(CF ₃) ₂	
949	i−C₃H ₇	Н	0	4-СН3	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
950	i-C ₃ H ₇	Н	0	4−CH ₃	2-C1-4-CF(CF ₃) ₂	
951	i−C₃H ₇	Н	0	4-CH₃	2-C1-4-(CF ₂) ₃ CF ₃	
952	i−C₃H ₇	Н	0	4−CH₃	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
953	i-C₃H₁	Н	0	4-СН₃	2-F-4-CF ₂ CF ₃	
954	i-C₃H₁	Н	0	4-CF ₃	2-CH ₃ -4-OCF ₃	
955	i-C₃H ₇	Н	0	4-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	

Table 10 (cont'd)

	_					Physical
No	R¹	R²	р	Xn	Ym	property
						m.p. ℃
956	i−C₃H ₇	Н	0	4−CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
957	i−C₃H ₇	Н	0	4−CF₃	2-CH ₃ -4-CF(CF ₃) ₂	
958	i-C ₃ H ₇	Н	0	4-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
959	i-C ₃ H ₇	Н	0	4-CF ₃	2-C1-4-CF(CF ₃) ₂	
960	i-C₃H₁	Н	0	4-CF ₃	2-C1-4-(CF ₂) ₃ CF ₃	
961	i-C ₃ H ₇	Н	0	4-CF ₃	2-C ₂ H ₅ -4-CF (CF ₃) ₂	
962	i−C₃H ₇	Н	0	4-CF ₃	2-F-4-CF ₂ CF ₃	
963	i−C₃H ₇	Н	0	4-CF₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
964	i−C₃H ₇	Н	0	4−CF₃	2-Br-4-CF ₂ CF ₃	
965	i−C₃H ₇	Н	0	5−CH₃	2-CH ₃ -4-0CF ₃	
966	i-C₃H,	Н	0	5-CH₃	2-CH ₃ -4-CF ₂ CF ₃	
967	i−C₃H ₇	Н	0	5-CH ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
968	i−C₃H ₇	Н	0	5−CH₃	2-CH ₃ -4-CF(CF ₃) ₂	
969	i−C₃H ₇	Н	0	6-СН3	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
970	i−C₃H ₇	Н	0	6-СН3	2-C1-4-CF(CF ₃) ₂	
971	i-C ₃ H ₇	Н	0	6−СН₃	2-C1-4-(CF ₂) ₃ CF ₃	
972	i-C ₃ H ₇	Н	0	6−СН₃	2-C ₂ H ₅ -4-CF (CF ₃) ₂	
973	i-C ₃ H ₇	Н	0	6-СН3	2-F-4-CF ₂ CF ₃	
974	i-C ₃ H ₇	Н	0	5-CF ₃	2-CH ₃ -4-OCF ₃	
975	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₃	
976	i-C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
977	i-C₃H₁	Н	0	5-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	

Table 10 (cont'd)

No	R¹	R²	р	Xn	Ym	Physical property m.p. °C
978	i−C₃H ₇	Н	0	5-CF ₃	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
979	i−C₃H ₇	Н	0	5-CF ₃	2-C1-4-CF(CF ₃) ₂	
980	i−C₃H ₇	Н	0	5-CF ₃	2-C1-4-(CF ₂) ₃ CF ₃	
981	i−C₃H ₇	Н	0	5-CF ₃	2-C ₂ H ₅ -4-CF ₂ CF ₃	
982	i−C₃H₁	Н	0	5-CF₃	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
983	i−C₃H ₇	Н	0	6-CF₃	2-F-4-CF ₂ CF ₃	
984	i−C₃H ₇	Н	0	6-CF ₃	2-Br-4-CF ₂ CF ₃	
985	i−C₃H ₇	Н	0	6-CF ₃	2-Br-4-CF(CF ₃) ₂	

Table 11
$$(R^1 = CH(CH_3)CH_2SCH_3, R^3 = H, Z^1 = Z^2 = O, Het = Q4, B^1 = B^2 = B^3 = B^4 = C)$$

No	R²	p	Xn	Ym	Physical property m.p. °C
986	Н	0	Н	2-CH ₃ -4-0CF ₃	-
987	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
988	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
989	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	108-110
990	Н	1	Н	2-CH ₃ -4-CF(CF ₃) ₂	paste
991	Н	0	Н	2-C1-4-CF(CF ₃) ₂	

Table 11 (cont'd)

No	R²	р	Xn	Ym	Physical property m.p. °C
992	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	
993	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
994	Н	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
995	Н	0	Н	2-F-4-CF ₂ CF ₃	
996	Н	0	Н	2-Br-4-CF ₂ CF ₃	
997	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
998	Н	0	4-Cl	2-CH ₃ -4-CF ₂ CF ₃	
999	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	237-239
1000	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1001	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
1002	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₃	
1003	Н	0	4-I	2-C1-4-CF(CF ₃) ₂	
1004	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1005	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	

Table 12 $(R^1 = C(CH_3)_2 CH_2 SCH_3, R^3 = H, Z^1 = Z^2 = O, He t = Q4, B^1 = B^2 = B^3 = B^4 = C)$

					Physical
No	R²	р	Xn	Ym	property
					m.p. ℃
1006	Н	0	Н	2-CH ₃ -4-0CF ₃	
1007	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1008	Н	0	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
1009	Н	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	paste
1010	Н	0	Н	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
1011	Н	0	Н	2-C1-4-CF(CF ₃) ₂	
1012	Н	0	Н	2-C1-4-(CF ₂) ₃ CF ₃	
1013	Н	0	Н	2-C ₂ H ₅ -4-CF ₂ CF ₃	
1014	Н	0	Н	2-C ₂ H ₅ -4-CF (CF ₃) ₂	
1015	Н	0	Н	2-F-4-CF ₂ CF ₃	
1016	Н	0	Н	2-Br-4-CF ₂ CF ₃	
1017	Н	0	Н	2-Br-4-CF(CF ₃) ₂	
1018	Н	0	4-C1	2-CH ₃ -4-CF ₂ CF ₃	
1019	Н	0	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1020	Н	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1021	Н	0	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
1022	Н	0	4-I	2-CH ₃ -4-CF ₂ CF ₃	
1023	Н	0	4-I	2-C1-4-CF(CF ₃) ₂	
1024	Н	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1025	Н	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	

Table 13 ($R^2 = R^3 = H$, $Z^1 = Z^2 = O$, Het = Q4, p = 0, $B^1 = B^2 = B^3 = B^4 = C$)

				Physical
No	R^1	Xn	Ym	
110	K	All	1111	property
1000	GIV (GIV.) GIV. GG. VI			m.p. ℃
1026	CH (CH ₃) CH ₂ SC ₂ H ₅	H	2-CH ₃ -4-0CF ₃	
1027	CH(CH₃)CH₂SC₂H₅	H	2-CH ₃ -4-CF ₂ CF ₃	
1028	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
1029	CH(CH ₃)CH ₂ SC ₂ H ₅	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1030	$CH(CH_3)CH_2SC_2H_5$	4-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
1031	$CH(CH_3)CH_2SC_2H_5$	4-I	2-C1-4-CF(CF ₃) ₂	
1032	$C(CH_3)_2CH_2SC_2H_5$	Н	2-CH ₃ -4-OCF ₃	
1033	$C(CH_3)_2CH_2SC_2H_5$	Н	2-CH ₃ -4-CF ₂ CF ₃	
1034	$C(CH_3)_2CH_2SC_2H_5$	Н	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
1035	$C(CH_3)_2CH_2SC_2H_5$	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1036	$C(CH_3)_2CH_2SC_2H_5$	5-Br	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
1037	$C(CH_3)_2CH_2SC_2H_5$	5-I	2-C1-4-CF(CF ₃) ₂	
1038	CH(CH ₃)CH ₂ NHAc	4-C1	2-CH ₃ -4-OCF ₃	
1039	CH(CH₃)CH₂NHAc	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
1040	CH(CH ₃)CH ₂ NHAc	5-I	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
1041	CH(CH ₃)CH ₂ NHAc	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
1042	C(CH ₃) ₂ CH ₂ NHAc	4-C1	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
1043	C(CH ₃) ₂ CH ₂ NHAc	5-I	2-C1-4-CF (CF ₃) ₂	
1044	CH(CH ₃)C ₂ H ₄ OCH ₃	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1045	$CH(CH_3)C_2H_4OCH_3$	4-I	2-CH ₃ -4-CF(CH ₃) ₂	
1046	$C(CH_3)_2C_2H_4OCH_3$	5-I	2-CH ₃ -4-CF(CF ₃) ₂	

Table 14
$$(Z^1 = Z^2 = 0, Het = Q4, B^1 = B^2 = B^3 = B^4 = C)$$

		T	T		ı		I
							Physical
	No	R¹	R ²	R³	Xn	Ym	property
							m.p. ℃
	1047	C ₂ H ₅	C ₂ H ₅	СН₃	Н	2-CH ₃ -4-OCF ₃	
	1048	C ₂ H ₅	C ₂ H ₅	СН₃	4-C1	2-CH ₃ -4-CF ₂ CF ₃	paste
	1049	C ₂ H ₅	C ₂ H ₅	СН₃	4-C1	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
	1050	C ₂ H ₅	C ₂ H ₅	СНз	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
	1051	C ₂ H ₅	C_2H_5	СНз	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
	1052	C ₂ H ₅	C ₂ H ₅	СНз	4-I	2-C1-4-CF (CF ₃) ₂	
	1053	C ₂ H ₅	C ₂ H ₅	C_2H_5	4-I	2-CH ₃ -4-OCF ₃	
	1054	C ₂ H ₅	C ₂ H ₅	C_2H_5	4-C1	2-CH ₃ -4-CF ₂ CF ₃	
	1055	C ₂ H ₅	C ₂ H ₅	C_2H_5	4-C1	2-CH ₃ -4-CF ₂ CF ₂ CF ₃	
	1056	C ₂ H ₅	C ₂ H ₅	C ₂ H ₅	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
	1057	C ₂ H ₅	C ₂ H ₅	C ₂ H ₅	4-I	2-CH ₃ -4-(CF ₂) ₃ CF ₃	
	1058	C ₂ H ₅	C ₂ H ₅	C_2H_5	4-I	2-C1-4-CF(CF ₃) ₂	
Į							

Table 15 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q5, B^1 = B^2 = B^3 = B^4 = C)$

					Physical
No	R¹	p	Xn	Ym	property
					m.p. ℃
1059	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1060	i−C₃H ₇	0	Н	2-CH ₃ -4-0CF ₃	
1061	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1062	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1063	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1064	i−C₃H ₇	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1065	i−C₃H ₇	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
1066	i−C₃H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1067	i−C₃H ₇	0	Н	2-C1-4-CF(CF ₂) ₂	
1068	i−C₃H ₇	0	Н	2-C1-4-CF ₂ CF ₃	
1069	i-C ₃ H ₇	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1070	i−C₃H ₇	0	Н	2-F-4-CF ₂ CF ₃	
1071	i-C ₃ H ₇	0	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
1072	$i-C_3H_7$	0	5-C1	2-CH ₃ -4-CF(CF ₃) ₂	,
1073	i-C ₃ H ₇	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
1074	i−C₃H ₇	0	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1075	i-C₃H ₇	0	5−СН₃	2-CH ₃ -4-CF ₂ CF ₃	
1076	i-C₃H₁	0	5-СН₃	2-CH ₃ -4-CF (CF ₃) ₂	
1077	i-C₃H ₇	0	5-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	

Table 15 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. ℃
1078	CH(CH ₃)CH ₂ SCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1079	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1080	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1081	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1082	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1083	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1084	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1085	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1086	$C(CH_3)_2CH_2NHAc$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1087	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1088	$CH(CH_3)C_2H_4OCH_3$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1089	$CH(CH_3)C_2H_4OCH_3$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 16 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q6, B^1 = B^2 = B^3 = B^4 = C)$

	1				···
No	R¹	р	Хn	Ym	Physical property
					m.p. ℃
1090	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1091	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₃	
1092	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1093	i−C₃H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1094	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1095	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	191-193
1096	i−C₃H ₇	0	Н	2-CH3-4-SCH2CF2CHF2	
1097	i−C₃H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1098	i−C₃H ₇	0	Н	2-C1-4-CF(CF ₂) ₂	
1099	i−C₃H ₇	0	Н	2-C1-4-CF ₂ CF ₃	
1100	i−C₃H ₇	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1101	i-C ₃ H ₇	0	Н	2-F-4-CF ₂ CF ₃	
1102	i−C₃H ₇	0	3-C1	2-CH ₃ -4-CF ₂ CF ₃	
1103	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1104	i-C ₃ H ₇	0	3-I	2-CH ₃ -4-CF ₂ CF ₃	
1105	i-C₃H ₇	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
1106	i−C₃H ₇	0	3-СН₃	2-CH ₃ -4-CF ₂ CF ₃	ļ
1107	i−C₃H ₇	0	6−СН₃	2-CH ₃ -4-CF(CF ₃) ₂	
1108	i−C₃H ₇	0	3-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	

Table 16 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. °C
1109	CH(CH ₃)CH ₂ SOCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1110	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	85-95
1111	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1112	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1113	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1114	C(CH ₃) ₂ CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1115	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1116	CH(CH₃)CH₂NHAc	0	Н	2 -CH $_3$ - 4 -CF(CF $_3$) $_2$	
1117	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1118	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2 -CH $_3$ - 4 -CF(CF $_3$) $_2$	
1119	$CH(CH_3)C_2H_4OCH_3$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	·
1120	$CH(CH_3)C_2H_4OCH_3$	0	Н	2 -CH ₃ - 4 -CF (CF ₃) $_2$	
1121	CH(CH₃)C₂H₄OCH₃	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
		:			

Table 17 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q7, B^1 = B^2 = B^3 = B^4 = C)$

	-1				Physical
No	R¹	р	Xn	n Ym	property
					m.p. °C
1122	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1123	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₃	
1124	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1125	i−C₃H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1126	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1127	i−C₃H ₇	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	204-206
1128	i-C ₃ H ₇	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
1129	i−C₃H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1130	$i-C_3H_7$	0	Н	2-C1-4-CF(CF ₂) ₂	
1131	i-C ₃ H ₇	0	Н	2-C1-4-CF ₂ CF ₃	
1132	i−C₃H₁	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1133	i−C₃H ₇	0	Н	2-F-4-CF ₂ CF ₃	,
1134	i−C₃H ₇	0	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
1135	i−C₃H ₇	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1136	i-C ₃ H ₇	0	5-I	2-CH ₃ -4-CF ₂ CF ₃	
1137	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
1138	i−C₃H ₇	0	5-CH₃	2-CH ₃ -4-CF ₂ CF ₃	
1139	i-C ₃ H ₇	0	6-СН ₃	2-CH ₃ -4-CF(CF ₃) ₂	
1140	i-C ₃ H ₇	0	5-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	

Table 17 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. °C
1141	CH(CH₃)CH₂SOCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1142	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1143	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1144	$CH(CH_3)CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1145	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1146	$C(CH_3)_2CH_2SOC_2H_5$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1147	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1148	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1149	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1150	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1151	CH(CH ₃)C ₂ H ₄ OCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1152	CH(CH₃)C₂H₄OCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 18 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q8, B^1 = B^2 = B^3 = B^4 = C)$

					Physical
No	R¹	р	Xn	Ym	property
					m.p. ℃
1153	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1154	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₃	
1155	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1156	i−C₃H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1157	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1158	i−C₃H ₇	0	Н	2-CH ₃ -4-CF (CF ₃) ₂	192-194
1159	i−C₃H₁	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
1160	i−C₃H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1161	i−C₃H ₇	0	Н	2-C1-4-CF(CF ₂) ₂	
1162	i−C₃H ₇	0	Н	2-C1-4-CF ₂ CF ₃	
1163	i-C ₃ H ₇	0	Н	2-C ₂ H ₅ -4-CF (CF ₃) ₂	
1164	i-C ₃ H ₇	0	Н	2-F-4-CF ₂ CF ₃	
1165	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF ₂ CF ₃	
1166	i−C₃H ₇	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1167	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF ₂ CF ₃	
1168	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
1169	i-C ₃ H ₇	0	2-СН3	2-CH ₃ -4-CF ₂ CF ₃	
1170	i−C₃H ₇	0	2-SCH ₃	2-CH ₃ -4-CF(CF ₃) ₂	204-206
1171	i-C ₃ H ₇	0	6−CF₃	2-CH ₃ -4-CF(CF ₃) ₂	
1172	CH(CH₃)CH₂	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SOCH ₃				

Table 18 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. °C
1173	CH(CH₃)CH₂SCH₃	0	2-SCH₃	2-CH ₃ -4-CF(CF ₃) ₂	168-170
1174	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1175	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1176	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1177	$C(CH_3)_2CH_2SOC_2H_5$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1178	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1179	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1180	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1181	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1182	CH(CH₃)C₂H₄OCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1183	CH(CH ₃)C ₂ H ₄ OCH ₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 19 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q9, B^1 = B^2 = B^3 = B^4 = C)$

	1	T		T	
No	R¹	p	Xn	Ym	Physical property
-					m.p. ℃
1184	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1185	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₃	
1186	i−C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1187	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1188	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1189	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	162-164
1190	i-C ₃ H ₇	0	Н	2-CH ₃ -4-SCH ₂ CF ₂ CHF ₂	
1191	i-C ₃ H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1192	i-C ₃ H ₇	0	Н	2-C1-4-CF(CF ₂) ₂	
1193	i-C ₃ H ₇	0	Н	2-C1-4-CF ₂ CF ₃	l.
1194	i-C₃H ₇	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1195	i-C ₃ H ₇	0	Н	2-F-4-CF ₂ CF ₃	
1196	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF ₂ CF ₃	
1197	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1198	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF ₂ CF ₃	
1199	i−C₃H ₇	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
1200	i-C ₃ H ₇	0	2-СН3	2-CH ₃ -4-CF ₂ CF ₃	
1201	i−C₃H ₇	0	2-SCH ₃	2-CH ₃ -4-CF(CF ₃) ₂	
1202	i-C ₃ H ₇	0	6-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	
1203	CH(CH ₃)CH ₂	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SOCH ₃				

Table 19 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. °C
1204	CH(CH₃)CH₂SCH₃	0	2-SCH ₃	2-CH ₃ -4-CF(CF ₃) ₂	129-131
1205	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1206	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1207	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1208	$C(CH_3)_2CH_2SOC_2H_5$	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1209	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1210	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1211	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	į
1212	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	1
1213	CH (CH ₃) C ₂ H ₄ OCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1214	CH(CH ₃)C ₂ H ₄ OCH ₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 20 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, q = 0, Het = Q10, B^1 = B^2 = B^3 = B^4 = C)$

	,	Γ	1		1
No	R ¹		v		Physical
NO	Т	P	Xn	Ym	property
					m.p. °C
1215	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1216	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₃	i
1217	i-C₃H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1218	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1219	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1220	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	153-155
1221	i−C₃H ₇	1	Н	2-CH ₃ -4-CF(CF ₃) ₂	paste
1222	i−C₃H ₇	0	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1223	i−C₃H ₇	0	Н	2-C1-4-CF(CF ₂) ₂	
1224	i−C₃H ₇	0	Н	2-C1-4-CF ₂ CF ₃	
1225	i−C₃H ₇	0	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1226	i−C₃H ₇	0	Н	2-F-4-CF ₂ CF ₃	
1227	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF ₂ CF ₃	
1228	i-C ₃ H ₇	0	6-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1229	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF ₂ CF ₃	
1230	i-C ₃ H ₇	0	6-I	2-CH ₃ -4-CF(CF ₃) ₂	
1231	i-C ₃ H ₇	0	2-СНз	2-CH ₃ -4-CF ₂ CF ₃	
1232	i-C₃H ₇	0	2-СН3	2-CH ₃ -4-CF(CF ₃) ₂	
1233	i-C₃H ₇	0	6-CF ₃	2-CH ₃ -4-CF(CF ₃) ₂	
1234	CH (CH ₃) CH ₂	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SCH₃				

Table 20 (cont'd)

No	R¹	р	Xn	Ym	Physical property m.p. °C
1235	CH(CH₃)CH₂SOCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1236	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1237	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1238	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1239	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1240	C(CH ₃) ₂ CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1241	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1242	CH(CH₃)CH₂NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1243	C(CH ₃)₂CH₂NHAc	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1244	C(CH ₃) ₂ CH ₂ NHAc	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1245	CH(CH₃)C₂H₄OCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1246	CH(CH ₃)C ₂ H ₄ OCH ₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 21 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q11, B^1 = B^2 = B^3 = B^4 = C)$

No	R¹	w	Xn	Ym	Physical property
					m.p. ℃
1247	i−C₃H₁	0	Н	2-CH ₃ -4-CF ₂ CF ₃	145-146
1248	i-C ₃ H ₇	0	4,5-(CH ₃) ₂	4-0CF ₃	148
1249	t-C ₄ H ₉	0	Н	2-CH ₃ -4-CF ₂ CF ₃	139-141
1250	i-C₃H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂	
				-CHF ₂	
1251	i−C₃H ₇	S	Н	2 -CH $_3$ - 4 -CF $_2$ CF $_3$	164-165
1252	i−C₃H ₇	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1253	i-C₃H₁	S	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
1254	i−C₃H ₇	S	Н	2-CH ₃ - 4 -SO ₂ CH ₂ CF ₂	
				-CHF ₂	
1255	i−C₃H ₇	S	4-I	2-CH ₃ -4-CF ₂ CF ₃	202-204
1256	t-C ₄ H ₉	S	Н	2-CH ₃ -4-CF ₂ CF ₃	178-180
1257	i−C₃H ₇	S	Н	$2-C_2H_5-4-CF(CF_3)_2$	
1258	i−C₃H ₇	NCH ₃	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1259	i−C₃H ₇	NCH₃	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
1250	i−C₃H₁	NCH₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1251	i−C₃H₁	NCH ₃	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1252	i−C₃H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1253	i−C₃H ₇	NPh	4-C1	2-CH ₃ -4-CF(CF ₃) ₂	
1254	i−C ₃ H ₇	NCH ₃	4-CH=CF	2-CH ₃ -4-CF(CF ₃) ₂	189-191
			-СН=СН-5		

Table 21 (cont'd)

No	R¹	w	Xn	Ym	Physical property
					m.p. ℃
1255	CH (CH ₃) CH ₂	NCH ₃	4-CH=CF	2-CH ₃ -4-CF(CF ₃) ₂	171-173
	−SCH ₃		-СН=СН-5		
1256	CH(CH₃)	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
	-C₂H₄OCH₃				
1257	CH (CH ₃) CH ₂	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SCH₃				
1258	$CH(CH_3)CH_2$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SOCH₃				;
1259	CH(CH₃)CH₂	S	Н	2-CH ₃ - 4 -CF(CF ₃) ₂	
	-SOC ₂ H ₅				
1260	$CH(CH_3)CH_2$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
	−SC ₂ H ₅				
1261	C(CH ₃) ₂ CH ₂	S	Н	2 -CH $_3$ -4-CF(CF $_3$) $_2$	
	-SC ₂ H ₅				
1262	$C(CH_3)_2CH_2$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
	-SOC ₂ H ₅				
1263	CH (CH ₃) CH ₂	S	Н	2 -CH $_3$ -4-CF(CF $_3$) $_2$	
	-NHAc				:
1264	C (CH ₃) ₂ CH ₂	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
	-NHAc				
1265	CH(CH ₃)	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
	-C ₂ H ₄ OCH ₃				

Table 22 $(Z^1 = Z^2 = 0, R^3 = H, Het = Q11, B^1 = B^2 = B^3 = B^4 = C)$

No	R¹	R ³	2	W	Xn	Ym	Physical property m.p. °C
1266 1267	C ₂ H ₅	C ₂ H		0	Н	2-CH ₃ -4-CF ₂ CF ₃	92-93
	C ₂ H ₅	C ₂ H		0	Н	2-CH ₃ -4-OCF ₃	81–82
1268	C ₂ H ₅	C ₂ H	5	S	Н	2-CH₃-4-CF₂CF₃	160-162
	[2-CH ₃ -	4-	Н	NCH ₃	4-CH=CF	2-CH ₃ -4-CF(CF ₃) ₂	256-258
1269	CF(CF ₃)	2]			-СН=СН-5		
	Ph						

Table 23
$$(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q12, B^1 = B^2 = B^3 = B^4 = C)$$

No	R¹	W	Xn	Ym	Physical property m.p. °C
1270	i−C₃H ₇	0	Н	2-CH ₃ -4-CF ₃	
1271	i−C₃H ₇	0	Н	4-0CF ₃	170
1272	i−C₃H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1273	i-C₃H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1274	i−C₃H ₇	S	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
1275	i−C₃H₁	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	140-145

Table 23 (cont'd)

			Ī .		
					Physical
No	R¹	W	Xn	Ym	property
					m.p. ℃
1276	i−C₃H ₇	S	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
1277	i−C₃H ₇	S	Н	2-CH ₃ -4-SO ₂ CH ₂	
				-CF ₂ CHF ₂	
1278	i−C₃H ₇	S	Н	2-C1-4-CF(CF ₂) ₂	
1279	i−C₃H ₇	S	Н	2-CH ₃ -5-CF ₂ CF ₃	125-130
1280	i−C₃H ₇	S	Н	2-CH ₃ -3-CF ₂ CF ₃	paste
1281	i-C ₃ H ₇	S	Н	2-F-4-CF ₂ CF ₃	
1282	i−C₃H ₇	NCH₃	2-I	2-CH ₃ -4-CF ₂ CF ₃	
1283	i−C₃H ₇	NCH ₃	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
1284	i-C ₃ H ₇	NCH ₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1285	i−C₃H ₇	NCH₃	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1286	i−C₃H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1287	i−C₃H ₇	NPh	2-I	2-CH ₃ -4-CF(CF ₃) ₂	
1288	i−C₃H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	: !
1289	CH(CH ₃)CH ₂ SCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1290	CH(CH ₃)CH ₂ SOCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1291	CH(CH ₃)CH ₂ SCH ₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1292	$CH(CH_3)CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1293	$CH(CH_3)CH_2SC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1294	$C(CH_3)_2CH_2SC_2H_5$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1295	$C(CH_3)_2CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 23 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1296	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1297	CH(CH ₃)CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1298	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1299	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1300	CH(CH₃)C₂H₄OCH₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1301	CH(CH₃)C₂H₄OCH₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 24 $(Z^1 = Z^2 = 0, R^3 = H, Het = Q13, B^1 = B^2 = B^3 = B^4 = C)$

				T		T	
	. T		_ 2				Physical
ľ	No	R¹	R ²	W	Xn	Ym	property
							m.p. ℃
13	302	i-C ₃ H ₇	Н	0	Н	2-CH ₃ -4-CF ₃	
13	303	i-C₃H₁	Н	0	4, 5-	4-0CF ₃	134
					(CH ₃) ₂		
13	304	i−C₃H₁	Н	0	Н	2-CH ₃ -4-0CF ₂ CHF ₂	
13	305	i-C₃H,	Н	S	Н	2-CH ₃ -4-OCH ₂ CF ₂	
						-CHF ₂	
13	806	i−C₃H ₇	Н	S	Н	2-CH ₃ -4-OCF ₃	139-141
13	07	i-C₃H ₇	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	159-161
13	808	i−C₃H₁	Н	S	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
13	09	i−C₃H ₇	Н	S	H	2-CH ₃ -4-CF(CF ₃) ₂	
13	10	i−C₃H ₇	Н	S	Н	2-C1-4-CF(CF ₂) ₂	
13	11	C ₂ H ₅	C_2H_5	S	Н	2-CH ₃ -4-CF ₂ CF ₃	120-122
13	12	C ₂ H ₅	C_2H_5	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
13	13	C ₂ H ₅	C_2H_5	S	Н	2-CH ₃ -4-OCF ₃	153-155
13	14	i-C ₃ H ₇	Н	NCH₃	4-I	2-CH ₃ -4-CF ₂ CF ₃	
13	15	i-C ₃ H ₇	Н	NPh	4-I	2-CH ₃ -4-CF(CF ₃) ₂	
13	16	i-C ₃ H ₇	Н	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
13	17	i-C ₃ H ₇	Н	NPh	Н	2-CH ₃ -4-CF(CF ₃) ₂	
13	18	i-C ₃ H ₇	Н	NСНз	Н	2-CH ₃ -4-CF ₂ CF ₃	
13	19	i-C ₃ H ₇	Н	NCH ₃	4-I	2-CH ₃ -4-CF (CF ₃) ₂	
<u></u>							

Table 24 (cont'd)

No	R¹	R²	W	Xn	Ym	Physical property m.p. °C
1320	i−C₃H ₇	Н	NCH₃	4-CH=CF -CH=CH-5	2-CH ₃ -4-CF (CF ₃) ₂	195-205
1321	CH(CH ₃)CH ₂ -SCH ₃ CH(CH ₃)CH ₂ -SOCH ₃	Н	S S	Н	2-CH ₃ -4-CF ₂ CF ₃ 2-CH ₃ -4-CF ₂ CF ₃	

Table 24 (cont'd)

No	R¹	R²	W	Xn	Ym	Physical property m.p. °C
1323	CH(CH₃)CH₂SCH₃	Н	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1324	CH(CH ₃)CH ₂ SOC ₂ H ₅	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1325	CH(CH₃)CH₂SC₂H₅	Н	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1326	$C(CH_3)_2CH_2SC_2H_5$	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1327	C(CH ₃) ₂ CH ₂ SOC ₂ H ₅	Н	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1328	CH(CH₃)CH₂NHAc	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1329	CH(CH₃)CH₂NHAc	Н	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1330	C(CH ₃) ₂ CH ₂ NHAc	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1331	C(CH ₃) ₂ CH ₂ NHAc	Н	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1332	CH(CH ₃)C ₂ H ₄ OCH ₃	Н	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1333	CH(CH ₃)C ₂ H ₄ OCH ₃	Н	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	

Table 25 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q14, B^1 = B^2 = B^3 = B^4 = C)$

	T		T T		T
No	R ¹	W	Xn	Ym	Physical property
					m.p. ℃
1334	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₃	
1335	i-C₃H₁	0	3-СН ₃	4-0CF ₃	137-138
1336	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1337	i-C ₃ H ₇	s	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1338	i-C ₃ H ₇	S	3-C1	2-CH ₃ -4-CF ₂ CF ₃	
1339	i-C ₃ H ₇	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1340	i-C ₃ H ₇	S	3-I	2-CH ₃ -4-CF(CF ₃) ₂	
1341	i-C ₃ H ₇	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1342	i-C ₃ H ₇	S	Н	2-C1-4-CF(CF ₂) ₂	
1343	i-C ₃ H ₇	S	Н	2-C1-4-CF ₂ CF ₃	
1344	i-C ₃ H ₇	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1345	i-C ₃ H ₇	S	Н	2-F-4-CF ₂ CF ₃	
1346	i−C₃H ₇	NCH ₃	3-I	2-CH ₃ -4-CF ₂ CF ₃	
1347	i-C ₃ H ₇	NCH ₃	3-I	2-CH ₃ -4-CF(CF ₃) ₂	
1348	i-C ₃ H ₇	NCH ₃	Н	2-CH ₃ -4-CF ₂ CF ₃	*
1349	i−C₃H ₇	NCH ₃	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1350	i−C₃H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1351	i-C₃H₁	NPh	3-I	2-CH ₃ -4-CF (CF ₃) ₂	
1352	i−C₃H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	

Table 25 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1352	CH(CH₃)CH₂SCH₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1353	CH(CH₃)CH₂SOCH₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1354	CH(CH₃)CH₂SCH₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1355	CH(CH ₃)CH ₂ SOC ₂ H ₅	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1356	CH(CH₃)CH₂SC₂H₅	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1357	$C(CH_3)_2CH_2SC_2H_5$	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1358	$C(CH_3)_2CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1359	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1360	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1361	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1362	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1363	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1364	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	

Table 26 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q15, B^1 = B^2 = B^3 = B^4 = C)$

	T	T	Т	T	· · · · · · · · · · · · · · · · · · ·
No	R ¹	W	Xn	Ym	Physical property
					m.p. °C
1365	i-C₃H₁	0	Н	2-CH ₃ -4-CF ₃	
1366	i-C₃H₁	0	5-СН3	4-0CF ₃	
1367	i-C₃H₁	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1368	i−C₃H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1369	i-C₃H₁	S	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
1370	i−C₃H ₇	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1371	i−C₃H ₇	S	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1372	i−C₃H ₇	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1373	i-C₃H₁	S	Н	2-C1-4-CF(CF ₂) ₂	
1374	i−C₃H ₇	S	Н	2-C1-4-CF ₂ CF ₃	
1375	i-C₃H₁	S	Н	2-C ₂ H ₅ -4-CF (CF ₃) ₂	
1376	i−C₃H ₇	S	Н	2-F-4-CF ₂ CF ₃	
1377	i−C₃H₁	NCH ₃	5-I	2-CH ₃ -4-CF ₂ CF ₃	
1378	i−C₃H₁	NCH ₃	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1379	i−C₃H ₇	NCH ₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1380	i−C₃H ₇	NCH ₃	5-Br	2-CH ₃ -4-CF(CF ₃) ₂	134-136
1381	i-C ₃ H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1382	i-C ₃ H ₇	NPh	5-Br	2-CH ₃ -4-CF(CF ₃) ₂	164-166
1383	i−C₃H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	

Table 26 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1384	CH(CH ₃)CH ₂ SCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1385	CH(CH₃)CH₂SOCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1386	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1387	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1388	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1389	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1390	C(CH ₃) ₂ CH ₂ SOC ₂ H ₅	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1391	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1392	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1393	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1394	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1395	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1396	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 27 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q16, B^1 = B^2 = B^3 = B^4 = C)$

No	R¹	W	Xn	Ym	Physical property m.p. °C
1397	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₃	ш. р. С
1398	i−C₃H ₇	0	5-CH ₃	4-0CF ₃	
1399	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1400	i-C ₃ H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1401	i−C₃H ₇	S	5-C1	2-CH ₃ -4-CF ₂ CF ₃	
1402	i−C₃H ₇	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1403	i-C₃H ₇	S	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1404	i-C ₃ H ₇	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1405	i-C₃H₁	S	Н	2-C1-4-CF(CF ₂) ₂	
1406	i-C ₃ H ₇	S	Н	2-C1-4-CF ₂ CF ₃	
1407	i-C ₃ H ₇	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1408	i−C₃H ₇	S	Н	2-F-4-CF ₂ CF ₃	
1409	i-C₃H ₇	NCH₃	5-I	2-CH ₃ -4-CF ₂ CF ₃	
1410	i-C₃H ₇	NCH₃	5-I	2-CH ₃ -4-CF(CF ₃) ₂	
1411	i−C₃H ₇	NCH₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1412	i−C₃H ₇	NCH₃	5-Br	2-CH ₃ -4-CF(CF ₃) ₂	165-175
1413	i−C₃H₁	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1414	i-C ₃ H ₇	NPh	5-Br	2-CH ₃ -4-CF(CF ₃) ₂	167-169
1415	i−C₃H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	
1416	CH(CH ₃)	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
	-CH ₂ SCH ₃				

Table 27 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1417	CH(CH ₃)CH ₂ SOCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1418	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1419	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1420	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	:
1421	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1422	C(CH ₃) ₂ CH ₂ SOC ₂ H ₅	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1422	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1423	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1424	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1425	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2 -CH $_3$ - 4 -CF(CF $_3$) $_2$	
1426	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1427	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 28 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q17, B^1 = B^2 = B^3 = B^4 = C)$

			T	·	T	
	No	R¹	W	Xn	Ym	Physical property
	1.400					m. p. ℃
	1428	i−C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₃	
	1429	i-C₃H ₇	0	3-CH ₃	4-0CF ₃	171-174
	1430	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
	1431	i-C ₃ H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
	1432	i-C ₃ H ₇	S	3-C1	2-CH ₃ -4-CF ₂ CF ₃	
	1433	i-C₃H ₇	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
	1434	i-C₃H ₇	S	3-I	2-CH ₃ -4-CF(CF ₃) ₂	
	1435	i−C₃H₁	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
	1436	i-C₃H ₇	S	Н	2-Cl-4-CF(CF ₂) ₂	
	1437	i−C₃H ₇	S	Н	2-C1-4-CF ₂ CF ₃	
	1438	i−C₃H ₇	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
	1439	i-C₃H₁	S	Н	2-F-4-CF ₂ CF ₃	
	1440	i−C₃H ₇	NCH ₃	3-I	2-CH ₃ -4-CF ₂ CF ₃	
	1441	i−C₃H ₇	NCH ₃	3-I	2-CH ₃ -4-CF(CF ₃) ₂	
	1442	i-C₃H₁	NCH ₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
	1443	i−C₃H₁	NCH ₃	3-Br	2-CH ₃ -4-CF(CF ₃) ₂	
	1444	i-C ₃ H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
	1445	i-C₃H₁	NPh	3-Br	2-CH ₃ -4-CF(CF ₃) ₂	
	1446	i−C₃H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	
L						

Table 28 (cont'd)

					Physical
No	R¹	W	Xn	Ym	property
					m.p. °C
1447	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1448	CH(CH₃)CH₂SOCH₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1449	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1450	CH(CH₃)CH₂SOC₂H₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1451	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1452	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1453	$C(CH_3)_2CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1454	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1455	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1456	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1457	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1458	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1459	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 29 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q18, B^1 = B^2 = B^3 = B^4 = C)$

	T	T			
No	R¹	W	Xn	Ym	Physical property m.p. °C
1460	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₃	
1461	i-C ₃ H ₇	0	Н	4-0CF ₃	
1462	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1463	i-C ₃ H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1464	i-C₃H ₇	S	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
1465	i−C₃H ₇	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1466	i−C₃H ₇	S	2-I	2-CH ₃ -4-CF (CF ₃) ₂	
1467	i-C ₃ H ₇	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1468	i-C ₃ H ₇	S	Н	2-C1-4-CF(CF ₂) ₂	
1469	i−C₃H₁	S	Н	2-C1-4-CF ₂ CF ₃	
1470	i-C₃H₁	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1471	i−C₃H ₇	S	Н	2-F-4-CF ₂ CF ₃	
1472	i−C₃H ₇	NCHF ₂	2-I	2-CH ₃ -4-CF ₂ CF ₃	
1473	i−C₃H ₇	NCHF ₂	Н	2-CH ₃ -4-CF(CF ₃) ₂	85-95
1474	i-C₃H₁	NCH₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1475	i−C₃H ₇	NCH ₃	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
1476	i−C ₃ H ₇	NPh	Н	2 -CH $_3$ -4-CF $_2$ CF $_3$	
1477	i-C ₃ H ₇	NPh	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
1478	i-C ₃ H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	

Table 29 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1479	CH(CH ₃)CH ₂ SCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1480	CH(CH ₃)CH ₂ SOCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1481	CH(CH₃)CH₂SCH₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1482	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1483	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1484	$C(CH_3)_2CH_2SC_2H_5$	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1485	$C(CH_3)_2CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1486	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1487	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1488	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1489	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1490	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	}
1491	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	

Table 30 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q19, B^1 = B^2 = B^3 = B^4 = C)$

			 -		
No	R ¹				Physical
NO	K	W	Xn	Ym	property
					m.p. ℃
1492	i-C ₃ H ₇	0	Н	2-CH ₃ -4-CF ₃	
1493	i-C ₃ H ₇	0	Н	4-0CF ₃	
1494	i-C ₃ H ₇	0	Н	2-CH ₃ -4-OCF ₂ CHF ₂	
1495	i-C ₃ H ₇	S	Н	2-CH ₃ -4-OCH ₂ CF ₂ CHF ₂	
1496	i-C ₃ H ₇	S	2-C1	2-CH ₃ -4-CF ₂ CF ₃	
1497	i-C ₃ H ₇	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1498	i-C ₃ H ₇	S	2-I	2-CH ₃ -4-CF (CF ₃) ₂	
1499	i−C₃H₁	S	Н	2-CH ₃ -4-SO ₂ CH ₂ CF ₂ CHF ₂	
1500	i-C₃H ₇	S	Н	2-Cl-4-CF(CF ₂) ₂	
1501	i−C₃H ₇	S	Н	2-C1-4-CF ₂ CF ₃	
1502	i-C ₃ H ₇	S	Н	2-C ₂ H ₅ -4-CF(CF ₃) ₂	
1503	i-C₃H₁	S	Н	2-F-4-CF ₂ CF ₃	
1504	i−C₃H ₇	NCHF ₂	2-I	2-CH ₃ -4-CF ₂ CF ₃	
1505	i−C₃H₁	NCHF ₂	Н	2-CH ₃ -4-CF(CF ₃) ₂	70-90
1506	i−C₃H₁	NCH ₃	Н	2-CH ₃ -4-CF ₂ CF ₃	
1507	i-C ₃ H ₇	NCH ₃	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
1508	i-C ₃ H ₇	NPh	Н	2-CH ₃ -4-CF ₂ CF ₃	
1509	i-C ₃ H ₇	NPh	2-Br	2-CH ₃ -4-CF(CF ₃) ₂	
1510	i-C ₃ H ₇	NPh	Н	2-C1-4-CF ₂ CF ₃	

Table 30 (cont'd)

No	R¹	W	Xn	Ym	Physical property m.p. °C
1511	CH (CH ₃) CH ₂ SCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1512	CH(CH ₃)CH ₂ SOCH ₃	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1513	CH(CH ₃)CH ₂ SCH ₃	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1514	CH(CH ₃)CH ₂ SOC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1515	CH(CH ₃)CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1516	C(CH ₃) ₂ CH ₂ SC ₂ H ₅	0	Н	2-CH ₃ -4-CF ₂ CF ₃	
1517	$C(CH_3)_2CH_2SOC_2H_5$	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1518	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1519	CH(CH₃)CH₂NHAc	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	
1520	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1521	C(CH ₃) ₂ CH ₂ NHAc	S	Н	2-CH ₃ -4-CF(CF ₃) ₂	
1522	CH(CH ₃)C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF ₂ CF ₃	
1523	CH (CH ₃) C ₂ H ₄ OCH ₃	S	Н	2-CH ₃ -4-CF (CF ₃) ₂	

Table 31 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, B^1 = B^2 = B^3 = B^4 = C)$

		Т		T	
No	R¹	W	Het	Ym	Physical property m.p. °C
1524	i-C ₃ H ₇	0	Q20	2-CH ₃ -4-CF ₂ CF ₃	
1525	i-C ₃ H ₇	0	Q20	2-CH ₃ -4-CF(CF ₃) ₂	
1526	i-C ₃ H ₇	S	Q20	2-CH ₃ -4-CF ₂ CF ₃	
1527	i-C ₃ H ₇	S	Q20	2-CH ₃ -4-CF (CF ₃) ₂	
1528	i-C ₃ H ₇	NCH₃	Q20	2-CH ₃ -4-CF ₂ CF ₃	
1529	i-C ₃ H ₇	NCH₃	Q20	2-CH ₃ -4-CF(CF ₃) ₂	
1530	i-C₃H₁	NPh	Q20	2-CH ₃ -4-CF ₂ CF ₃	
1531	i−C₃H ₇	NPh	Q20	2-CH ₃ -4-CF(CF ₃) ₂	i
1532	i-C₃H₁	0	Q21	2-CH ₃ -4-CF ₂ CF ₃	
1533	i-C ₃ H ₇	0	Q21	2-CH ₃ -4-CF(CF ₃) ₂	
1534	i-C₃H₁	S	Q21	2-CH ₃ -4-CF ₂ CF ₃	
1535	i-C ₃ H ₇	S	Q21	2-CH ₃ -4-CF(CF ₃) ₂	
1536	i-C₃H₁	NCH ₃	Q21	2-CH ₃ -4-CF ₂ CF ₃	
1537	i−C₃H ₇	NCH ₃	Q21	2-CH ₃ -4-CF(CF ₃) ₂	
1538	i-C₃H₁	NPh	Q21	2-CH ₃ -4-CF ₂ CF ₃	
1539	i-C ₃ H ₇	NPh	Q21	2-CH ₃ -4-CF(CF ₃) ₂	
1540	i-C ₃ H ₇	0	Q22	2-CH ₃ -4-CF ₂ CF ₃	
1541	i-C₃H ₇	0	Q22	2-CH ₃ -4-CF(CF ₃) ₂	
1542	i−C ₃ H ₇	S	Q22	2-CH ₃ -4-CF ₂ CF ₃	
1543	i-C ₃ H ₇	S	Q22	2-CH ₃ -4-CF(CF ₃) ₂	
1544	i-C ₃ H ₇	NCH ₃	Q22	2-CH ₃ -4-CF ₂ CF ₃	

Table 31 (cont'd)

No	R¹	W	Het	Ym	Physical property m.p. °C
1545	i−C₃H₁	NCH₃	Q22	2-CH ₃ -4-CF(CF ₃) ₂	
1546	i−C₃H ₇	NPh	Q22	2-CH ₃ -4-CF ₂ CF ₃	
1547	i-C ₃ H ₇	NPh	Q22	2-CH ₃ -4-CF(CF ₃) ₂	

Table 32
$$(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q2, p = 0, B^1 = B^4 = C)$$

No	R¹	Xn	B ²	B ³	Ym	Physical property m.p. °C
1548	i-C ₃ H ₇	Н	N	С	2-CH ₃ -4-OCH(CF ₃) ₂	259-260
1549	C(CH ₃) ₂ CH ₂ SCH ₃	Н	N	С	2-CH ₃ -4-OCH(CF ₃) ₂	202-203
1550	C(CH ₃) ₂ CH ₂ SCH ₃	5-C1	N	С	2-CH ₃ -4-CF(CF ₃) ₂	
1551	i−C₃H ₇	Н	N	N	2 -CH $_3$ -4-CF(CF $_3$) $_2$	
1552	i−C₃H ₇	5-C1	С	N	2-CH ₃ -4-CF ₂ CF ₃	
1553	i−C₃H ₇	5-I	С	N	4-0CH(CF ₃) ₂	
1554	CH(CH₃)CH₂SCH₃	Н	N	N	4-0CHF ₂	
1555	C(CH ₃) ₂ CH ₂	Н	N	С	2-CH ₃ -4-OCF ₃	
	−SOCH ₃					
1556	CH(CH₃)CH(CH₃)	Н	N	С	2-CH ₃ -4-CF ₂ CF ₃	
	-SCH ₃					

Table 33 $(Z^1 = Z^2 = 0, R^2 = R^3 = H, Het = Q3, p = 0, B^1 = B^4 = C)$

No	R¹	Xn	B²	B³	Ym	Physical property m.p. °C
1557	i-C ₃ H ₇	Н	N	С	2-CH ₃ -4-OCH(CF ₃) ₂	
1558	C(CH ₃) ₂ CH ₂ SCH ₃	Н	N	С	2-CH ₃ -4-OCH(CF ₃) ₂	
1559	C(CH ₃) ₂ CH ₂ SCH ₃	2-C1	N	С	2-CH ₃ -4-CF(CF ₃) ₂	
1560	i−C₃H ₇	Н	N	N	2-CH ₃ -4-CF(CF ₃) ₂	
1561	i−C₃H ₇	2-C1	С	N	2-CH ₃ -4-CF ₂ CF ₃	
1562	i−C₃H ₇	2-I	С	N	4-0CH(CF ₃) ₂	
1563	CH(CH₃)CH₂SCH₃	Н	N	N	4-0CHF ₂	
1564	C (CH ₃) ₂ CH ₂	Н	N	С	2-CH ₃ -4-OCF ₃	
	-SOCH ₃	ı				
1565	СН (СНз) СН (СНз)	Н	N	С	2-CH ₃ -4-CF ₂ CF ₃	
	−SCH₃					

In the Tables 1 to 33, "Ac" means acetyl group, "Ph" means phenyl group and "c-" means alicyclic hydrocarbon group.

The agricultural and horticultural insecticides containing the heterocyclic dicarboxylic acid diamide derivative of formula (I) of the present invention as an active ingredient are suitable for 5 controlling various insect pests such as agrohorticultural insect pests, stored grain insect pests, sanitary insect pests, nematodes, etc., which are injurious to paddy rice, fruit trees, vegetables, other crops, flowers, ornamental plants, etc. 10 have a marked insecticidal effect, for example, on LEPIDOPTERA including summer fruit tortrix (Adoxophes orana fasciata), smaller tea tortrix (Adoxophyes sp.), Manchurian fruit moth (Grapholita inopinata), oriental fruit moth (Grapholita molesta), soybean pod border 15 (Leguminovora glycinivorella), mulberry leafroller (Olethreutes mori), tea leafroller (Caloptilia thevivora), Caloptilia sp. (Caloptilia zachrysa), apple leafminer (Phyllonorycter ringoniella), pear barkminer (Spulerrina astaurota), common white (Piers rapae 20 crucivora), tobacco budworm (Heliothis sp.), codling moth (Laspey resia pomonella), diamondback moth (Plutella xylostella), apple fruit moth (Argyresthia conjugella), peach fruit moth (Carposina niponensis), rice stem borer (Chilo suppressalis), rice leafroller (Cnaphalocrocis medinalis), tobacco moth (Ephestia 25 elutella), mulberry pyralid (Glyphodes pyloalis), yellow rice borer (Scirpophaga incertulas), rice

skipper (Parnara guttata), rice armyworm (Pseudaletia

separata), pink borer (Sesamia inferens), common
cutworm (Spodoptera litura), beet armyworm (Spodoptera
exigua), etc.; HEMIPTERA including aster leafhopper
(Macrosteles fascifrons), green rice leafhopper

- (Nephotettix cincticepts), brown rice planthopper (Nilaparvata lugens), whitebacked rice planthopper (Sogatella furcifera), citrus psylla (Diaphorina citri), grape whitefly (Aleurolibus taonabae), sweetpotato whitefly (Bemisia tabaci), greenhouse
- whitefly (Trialeurodes vaporariorum), turnup aphid
 (Lipaphis erysimi), green peach aphid (Myzus persicae),
 Indian wax scale (Ceroplastes ceriferus), cottony
 citrus scale (Pulvinaria aurantii), camphor scale
 (Pseudaonidia duplex), san Jose scale (Comstockaspis
- perniciosa), arrowhead scale (Unapsis yanonensis),
 etc.; TYLENCHIDA including soybean beetle (Anomala
 rufocuprea), Japanese beetle (Popillia japonica),
 tobacco beetle (Lasioderma serricorne), powderpost
 beetle (Lyctus brunneus), twenty-eight-spotted ladybird
- (Callosobruchus chinensis), vegetable weevil
 (Listroderes costirostris), maize weevil (Sitophilus zeamais), boll weevil (Anthonomus gradis gradis), rice water weevil (Lissorhoptrus oryzophilus), cucurbit leaf
- 25 beetle (<u>Aulacophora femoralis</u>), rice leaf beetle

 (<u>Oulema oryzae</u>), striped flea beetle (<u>Phyllotreta</u>

 <u>striolata</u>), pine shoot beetle (<u>Tomicus piniperda</u>),

 Colorado potato beetle (<u>Leptinotarsa decemlineata</u>),

Mexican bean beetle (Epilachna varivestis), corn rootworm (Diabrotica sp.), etc.; DIPTERA including (Dacus(Zeugodacus) cucurbitae), oriental fruit fly (Dacus(Bactrocera) dorsalis), rice leafminer (Agnomyza oryzae), onion maggot (Delia antiqua), seedcorn maggot (Delia platura), soybean pod gall midge (Asphondylia sp.), muscid fly (Musca domestica), house mosquito (Culex pipiens pipiens), etc.; and TYLENCHIDA including coffee root-lesion nematode (Pratylenchus coffeae), potato cyst nematode (Globodera rostochiensis), root-knot nematode (Meloidogyne sp.), citrus nematode (Tylenchulus semipenetrans), Aphelenchus sp. (Aphelenchus avenae), chrysanthemum foliar

15 The agricultural and horticultural insecticide containing the heterocyclic dicarboxylic acid diamide derivative represented by formula (I) of the present invention has a marked controlling effect on the above-exemplified insect pests, sanitary pests 20 and/or nematodes, which are injurious to paddy field crops, upland crops, fruit trees, vegetables and other crops, flowers and ornament plants, and the like. Therefore, the desired effect of the agrohorticultural insecticide of the present invention can be exhibited by applying the insecticide to the paddy field water, stalks and leaves or soil of paddy field, upland field, fruit trees, vegetables, other crops or flowers and ornament plants at a season at which the insect pests,

(Aphelenchoides ritzemabosi), etc.

sanitary pests or nematodes are expected to appear, before their appearance or at the time when their appearance is confirmed.

In general, the agricultural and

5 horticultural insecticide of the present invention is
used after being prepared into conveniently usable
forms according to ordinary manner for preparation of
agrochemicals.

That is, the heterocyclic dicarboxylic acid

diamide derivative of formula (I) and an appropriate

carrier are blended optionally together with an

adjuvant in a proper proportion and prepared into a

suitable preparation form such as suspension,

emulsifiable concentrate, soluble concentrate, wettable

powder, granules, dust or tablets through dissolution,

separation, suspension, mixing, impregnation, adsorption or sticking.

The inert carrier used in the present invention may be either solid or liquid. As the solid carrier, soybean flour, cereal flour, wood flour, bark flour, saw dust, powdered tobacco stalks, powdered walnut shells, bran, powdered cellulose, extraction residues of vegetables, powdered synthetic polymers or resins, clay (e.g. kaolin, bentonite and acid clay), talc (e.g. talc and pyrophyllite), silica materials (e.g. diatomaceous earth, siliceous sand, mica, white carbon, i.e. synthetic high-dispersion silicic acid, also called finely divided hydrated silica or hydrated

silicic acid, some of the commercially available products contain calcium silicate as the major component), activated carbon, powdered sulfur, pumice, calcined diatomaceous earth, ground brick, fly ash, sand, calcium carbonate, calcium phosphate and other inorganic or mineral powders, chemical fertilizers such as ammonium sulfate, ammonium phosphate, ammonium nitrate, urea, ammonium chloride and the like, and compost. These carriers may be used either alone or as a mixture of two or more carriers.

The liquid carrier is that which itself has a solubility or which is without such solubility but is capable of dispersing an active ingredient with the aid of an adjuvant. The following are typical examples of 15 the liquid carrier and can be used alone or as a mixture thereof. Water; alcohols such as methanol, ethanol, isopropanol, butanol and ethylene glycol; ketones such as acetone, methyl ethyl ketone, methyl isobutyl ketone, diisobutyl ketone and cyclohexanone; 20 ethers such as ethyl ether, dioxane, cellosolve, dipropyl ether and tetrahydrofuran; aliphatic hydrocarbons such as kerosene and mineral oil; aromatic hydrocarbons such as benzene, toluene, xylene, solvent naphtha and alkylnaphthalene; halogenated hydrocarbons 25 such as dichlorethane, chloroform, carbon tetrachloride and chlorobenzene; esters such as ethyl acetate, diisopropyl phthalate, dibutyl phthalate and dioctyl phthalate; amides such as dimethylformamide,

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diethylformamide and dimethylacetamide; nitriles such as acetonitrile; and dimethyl sulfoxide.

The following are typical examples of the adjuvant, which are used depending upon purposes and used alone or in combination of two or more adjuvants in some cases, or need not to be used at all.

To emulsify, disperse, dissolve and/or wet an active ingredient, a surfactant is used. As the surfactant, there can be exemplified polyoxyethylene alkyl ethers, polyoxyethylene alkylaryl ethers, polyoxyethylene higher fatty acid esters, polyoxyethylene resinates, polyoxyethylene sorbitan monolaurate, polyoxyethylene sorbitan monolaurate, polyoxyethylene sorbitan monooleate, alkylarylsulfonates, naphthalene-sulfonic acid condensation products, ligninsulfonates and higher alcohol sulfate esters.

Further, to stabilize the dispersion of an active ingredient, tackify it and/or bind it, there may be used adjuvants such as casein, gelatin, starch, methyl cellulose, carboxymethyl cellulose, gum arabic, polyvinyl alcohols, turpentine, bran oil, bentonite and ligninsulfonates.

To improve the flowability of a solid product, there may be used adjuvants such as waxes, stearates and alkyl phosphates.

Adjuvants such as naphthalenesulfonic acid condensation products and polycondensates of phosphates may be used as a peptizer for dispersible products.

Adjuvants such as silicone oil may also be used as a defoaming agent.

The content of the active ingredient may be varied according to the need. In dusts or granules, the suitable content thereof is from 0.01 to 50% by weight. In emulsifiable concentrate and flowable wettable powder, too, the suitable content is from 0.01 to 50% by weight.

insecticide of the present invention is used to control a variety of insect pests in the following manner.

That is, it is applied to a crop on which the insect pests are expected to appear or a site where appearance of the insect pests is undesirable, as it is or after being properly diluted with or suspended in water or the like, in an amount effective for control of the insect pests.

The applying dosage of the agrihorticultural insecticide of the present invention is varied

20 depending upon various factors such as a purpose, insect pests to be controlled, a growth state of a plant, tendency of insect pests appearance, weather, environmental conditions, a preparation form, an application method, an application site and an

25 application time. It may be properly chosen in a range of 0.1 g to 10 kg (in terms of active ingredient compound) per 10 are depending upon purposes.

The agricultural and horticultural

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insecticide of the present invention may be used in admixture with other agricultural and horticultural disease or pest controllers in order to expand both spectrum of controllable diseases and insect pest species and the period of time when effective applications are possible or to reduce the dosage.

EXAMPLES

Typical examples of the present invention are described below, but they should not be construed as limiting the scope of the invention.

Example 1

Methylphenyl]pyridine-3,4-dicarboximide

In 10 ml of THF were dissolved 1.50 g of

pyridine-3,4-dicarboxylic acid anhydride and 2.75 g of

4-(heptafluoro-2-propyl)-2-methylaniline, and the
reaction was carried out at room temperature for 3
hours. After completion of the reaction, the solvent
was distilled off under reduced pressure and 20 ml of

trifluoroacetic anhydride was added to the resulting
residue, and the reaction was carried out with
refluxing for 3 hours. After completion of the

25 methylphenyl]pyridine-3,4-dicarboximide quantitatively.

reaction, the solvent was distilled off under reduced

pressure to obtain crude N-[4-(heptafluoro-2-propyl)-2-

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(1-2). Production of 3-[4-(heptafluoro-2-propyl)-2methylphenyl]aminocarbonyl-4-pyridinecarboxylic
acid-2-propylamide (compound No. 230) and 4-[4(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-3-pyridinecarboxylic acid-2propylamide (compound No. 512)

In 10 ml of dioxane was dissolved 4.1 g of N[4-(heptafluoro-2-propyl)-2-methylphenyl]pyridine-3,4dicarboximide, followed by adding thereto 0.8 g of

10 isopropylamine, and the reaction was carried out at
room temperature for 8 hours. After completion of the
reaction, the solvent was distilled off under reduced
pressure and the resulting residue was purified by a
silica gel column chromatography using a 2 : 1 mixed

15 solvent of hexane and ethyl acetate as an eluent, to
obtain 2.1 g of (compound No. 230) and 1.8 g of
(compound No. 512) as white crystals.

Physical property:

compound No. 230 m.p. 234-236°C. Yield 45%.

compound No. 512 m.p. 206-208°C. Yield 39%.

Example 2

(2-1). Production of 5-bromo-3-[4-(heptafluoro-2propyl)-2-methylphenyl]aminocarbonyl-1-phenyl4-pyrazolecarboxylic acid 2-propylamide
(compound No. 1382) and 5-bromo-4-[4(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-1-phenyl-3-pyrazolecarboxylic

acid 2-propylamide (compound No. 1414)

In 10 ml of thionyl chloride was dissolved 500 mg of 5-bromo-1-phenyl-3,4-pyrazoledicarboxylic acid, and the reaction was carried out at reflux temperature for 2 hours. After completion of the reaction, the thionyl chloride was distilled off under reduced pressure to obtain a crude acid chloride. This compound was dissolved in 2 ml of THF and the resulting solution was added dropwise at 0°C to a solution

- prepared by dissolving 420 mg of heptafluoro-2-propyl-2-methylaniline and 410 mg of triethylamine in 10 ml of THF. After completion of the dropwise addition, 470 mg of isopropylamine was added thereto at 0°C and the reaction was carried out at room temperature for 2
- 15 hours. After completion of the reaction, the triethylamine hydrochloride was filtered off and the mother liquor was concentrated. The resulting residue was purified by a silica gel column chromatography using ethyl acetate/n-hexane as an eluent, to obtain
- 360 mg of 5-bromo-3-[4-(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-1-phenyl-4-pyrazole-carboxylic acid 2-propylamide (compound No. 1382) and 360 mg of 5-bromo-4-[4-(heptafluoro-2-propyl)-2-methylphenyl]-aminocarbonyl-1-phenyl-3-pyrazole-
- 25 carboxylic acid 2-propylamide (compound No. 1414) as white crystals.

Physical property:

compound No. 1381 m.p. $164-166^{\circ}$ C. Yield 36%.

compound No. 1412 m.p. 167-169°C. Yield 36%.

Example 3

5

- (3-1). Production of 2-chloro-4-[4-(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-3-pyridinecarboxylic acid
- In 20 ml of anhydrous THF was dissolved 2.78 g of diisopropylamine, and 18 ml of a solution of n-butyllithium in hexane (1.53 M) was added dropwise thereto at -78% under an argon atmosphere. The
- resulting mixture was stirred at -78℃ for 1 hour, and then a solution of 5.17 g of 2-chloro-4pyridinecarboxylic acid 4-(heptafluoro-2-propyl)-2methylanilide in 100 ml of THF was added dropwise
 thereto at -78℃. After completion of the dropwise
- 15 addition, the stirring was continued at -78℃ for 2 hours and carbon dioxide gas was bubbled thereinto for 1 hour. Then, the resulting mixture was warmed to room temperature and 200 ml of 1N hydrochloric acid was added thereto, followed by extraction with ethyl
- acetate. The organic layer was dried over anhydrous magnesium sulfate and then distilled under reduced pressure to remove the solvent, whereby 4.70 g (yield: 82%) of the crude desired compound was obtained as an amorphous solid. This compound was used in a
- 25 subsequent reaction without purification.

- (3-2). Production of 2-chloro-4-[4-(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-3-pyridinecarboxylic acid 2-propylamide (compound No. 524)
- In 10 ml of t-butyl methyl ether was dissolved 500 mg of 2-chloro-4-[4-(heptafluoro-2-propyl)-2-methylphenyl]aminocarbonyl-3-pyridine-carboxylic acid, followed by adding thereto 340 mg of trifluoroacetic anhydride, and the resulting mixture
- was stirred at room temperature for 2 hours. After the disappearance of the starting material was confirmed by TLC, 330 mg of isopropylamine was added to the mixture, followed by stirring at room temperature for another 2 hours. After completion of the reaction, ethyl acetate
- was added and the reaction solution was washed successively with water, a saturated aqueous sodium hydrogencarbonate solution and a saturated aqueous sodium chloride solution, and dried over anhydrous magnesium sulfate. The solvent was distilled off under
- reduced pressure and the resulting residue was purified by a silica gel column chromatography using ethyl acetate/n-hexane as an eluent, to obtain 460 mg of the desired compound as white crystals.

Physical property: m.p. 275-277℃. Yield 84%.

25 Example 4

(4-1). Production of N-[4-(heptafluoro-2-propyl)-2-methylphenyl]pyridine-2,3-dicarboximide-1-oxide

In 25 ml of chloroform was dissolved 3.1 g of N-[4-(heptafluoro-2-propyl)-2-methylphenyl]pyridine-2,3-dicarboximide, and 5.0 g of m-chloroperbenzoic acid was added thereto at room temperature. After the resulting mixture was stirred at room temperature for 3 hours, a saturated aqueous sodium hydrogencarbonate solution was added to the reaction solution, followed by extraction with ethyl acetate. The organic layer was washed successively with water, a saturated aqueous 10 sodium hydrogencarbonate solution and a saturated aqueous sodium chloride solution, and dried over anhydrous magnesium sulfate. The solvent was distilled off under reduced pressure and the resulting residue was purified by a silica gel column chromatography 15 using ethyl acetate/n-hexane as an eluent, to obtain 820 mg (yield: 84%) of the desired compound.

- (4-2). Production of 3-[4-(heptafluoro-2-propy1)-2-methylphenyl]aminocarbonyl-2-(2-propyl)aminocarbonylpyridine-N-oxide (compound No. 804)
- In 10 ml of THF was dissolved 400 mg of N-[4-(heptafluoro-2-propyl)-2-methylphenyl]pyridine-2,3-dicarboximide-1-oxide, followed by adding thereto 200 mg of isopropylamine, and the resulting mixture was stirred at room temperature for 5 hours. After

 25 completion of the reaction, the solvent was distilled off under reduced pressure and the resulting residue was purified by a silica gel column chromatography

using ethyl acetate/n-hexane as an eluent, to obtain 290 mg of the desired compound as white crystals.

Physical property: m.p. $108-110^{\circ}$ C. Yield 63%.

Typical formulation examples and test

5 examples of the present invention are described below but they should not be construed as limiting the scope of the invention.

In the formulation examples, parts are all by weight.

10 Formulation Example 1

Each compound listed in Tables 50 parts

1 to 33

Xylene 40 parts

Mixture of polyoxyethylene 10 parts

nonylphenyl ether and calcium alkylbenzenesulfonate

An emulsifiable concentrate was prepared by mixing uniformly the above ingredients to effect dissolution.

20 Formulation Example 2

Each compound listed in Tables 3 parts

1 to 33

Clay powder 82 parts

Diatomaceous earth powder 15 parts

A dust was prepared by mixing uniformly and

10

15

grinding the above ingredients.

Formulation Example 3

Each compound listed in Tables 5 parts

1 to 33

5 Mixed powder of bentonite and clay 90 parts
Calcium lignin sulfonate 5 parts

Granules were prepared by mixing the above ingredients uniformly, and kneading the resulting mixture together with a suitable amount of water, followed by granulation and drying.

Formulation Example 4

Each compound listed in Tables 20 parts

1 to 33

Kaolin and synthetic high— 75 parts
dispersion silicic acid

Mixture of polyoxyethylene 5 parts
nonylphenyl ether and calcium

A wettable powder was prepared by mixing 20 uniformly and grinding the above ingredients.

alkylbenzenesulfonate

Test Example 1

Insecticidal effect on diamondback moth (Plutella
xylostella)

Adult diamondback moths were released and 25 allowed to oviposit on a Chinese cabbage seedling. Two

days after the release, the seedling having eggs
deposited thereon was immersed for about 30 seconds in
a liquid chemical prepared by diluting a preparation
containing each of the compounds listed in Tables 1 to

33 as an active ingredient to adjust the concentration
to 500 ppm. After air-drying, the seedling was allowed
to stand in a room thermostated at 25°C. Six days after
the immersion in the liquid chemical, the hatched
insects were counted and the mortality was calculated

10 according to the following equation. The test was
carried out with three replications of 10 insects.

As a result, the following compounds were found to have a corrected mortality of 90% or more: compound Nos. 13, 18, 32, 54, 55, 57, 127, 136, 230, 242, 258, 464, 484, 512, 524, 737, 785, 794, 795, 804, 805, 821, 989, 990, 1009, 1095, 1110, 1127, 1158, 1189, 1204, 1220, 1221, 1247, 1249, 1251, 1255, 1267, 1269, 1271, 1275, 1303, 1306, 1313, 1320, 1414, 1429, 1473, 1505, 1548 and 1540.

Test Example 2

Insecticidal effect on smaller tea tortrix
(Adoxophyes sp.)

A tea leaf was immersed for about 30 seconds in a liquid chemical prepared by diluting a preparation containing each of the compounds listed in Tables 1 to 33 as an active ingredient to adjust the concentration to 500 ppm. After air-drying, the leaf was placed in a plastic Petri dish with a diameter of 9 cm and

- inoculated with larvae of smaller tea tortrix, and the Petri dish was allowed to stand in a room thermostated at 25℃ and having a humidity of 70%. Eight days after the inoculation, the dead and alive were counted and the mortality was calculated according to the following
- 15 equation. The test was carried out with three replications of 10 insects.

As a result, the following compounds were found to have a corrected mortality of 90% or more: compound Nos. 13, 18, 55, 57, 127, 136, 230, 464, 484, 512, 524, 737, 794, 795, 805, 821, 989, 1009, 1048, 1095, 1127, 1189, 1204, 1220, 1247, 1249, 1251, 1255,

1303, 1313, 1473, 1505, 1548 and 1549.

As described above, the agricultural and horticultural insecticides containing the heterocyclic dicarboxylic acid diamide derivative of the general formula (I) of the present invention as an active ingredient have an excellent controlling effect on insect pests such as diamondback moth, common cutworm, etc.

CLAIMS

1. A heterocyclic dicarboxylic acid diamide derivative represented by the general formula (I):

$$\begin{array}{c|c} Xn & Z^1 \\ \hline NR^1R^2 \\ \hline NR^2 \\ \hline NR^1R^2 \\ \hline$$

{wherein R^1 , R^2 and R^3 , which may be the same or different, are hydrogen atoms, (C_3-C_6) cycloalkyl groups, halo (C_3-C_6) cycloalkyl groups or $-A^1-(R^4)$ r (wherein A^1 is a (C_1-C_8) alkylene group, a (C_3-C_6) alkenylene group or a (C_3-C_6) alkynylene group, R^4 , which may be the same or different, are hydrogen atoms; halogen atoms; cyano groups; nitro groups; halo (C_1-C_6) alkyl groups; $(C_3 C_6$) cycloalkyl groups; halo (C_3-C_6) cycloalkyl groups; (C_1-C_6) C_6) alkoxycarbonyl groups; di(C_1 - C_6) alkoxyphosphoryl groups whose (C_1-C_6) alkoxy groups may be the same or different; $di(C_1-C_6)$ alkoxythiophosphoryl groups whose (C_1-C_6) alkoxy groups may be the same or different; diphenylphosphino groups; diphenylphosphono groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups,

halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^2-R^5$ (wherein A^2 is -O-, -S-, -SO-, -SO₂-, -N(R^6)- (wherein R^6 is a hydrogen atom; a (C_1-C_6) alkylcarbonyl group; a halo(C_1-C_6)alkylcarbonyl group; a (C_1-C_6)alkoxycarbonyl group; a phenylcarbonyl group; a substituted phenylcarbonyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) C_6) alkylsulfonyl groups; a phenyl (C_1-C_4) alkoxycarbonyl group; a substituted phenyl(C₁-C₄)alkoxycarbonyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6)

 C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a (C_1-C_6) alkylsulfonyl group; or a halo (C_1-C_6) C_6) alkylsulfonyl group), -C(=0) - or $-C(=NOR^7)$ - (wherein R^7 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo (C_1-C_6) C_6) alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) C_6) alkenyl group; a (C_3-C_6) alkynyl group; a cyclo (C_3-C_6) C_6) alkyl group; a phenyl (C_1-C_4) alkyl group; or a substituted phenyl(C_1-C_4)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1 - C_6) alkylthio groups, (C_1 - C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, $(C_1 C_6$) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and R^5 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo(C_1-C_6) alkyl group; a (C_3-C_6) alkenyl group; a halo (C_3-C_6) alkenyl group; a (C_3-C_6) alkynyl group; a halo(C_3-C_6) alkynyl group; a (C_3-C_6) cycloalkyl group; a $\label{eq:condition} \verb|halo(C_3-C_6)| \ cycloalkyl group; \ a \ (C_1-C_6) \ alkoxy(C_1-C_6) \ alkyl$ group; a (C_1-C_6) alkylthio (C_1-C_6) alkyl group; a formyl group; a (C_1-C_6) alkylcarbonyl group; a halo $(C_1 C_6$) alkylcarbonyl group; a (C_1-C_6) alkoxycarbonyl group; a $mono(C_1-C_6)$ alkylaminocarbonyl group; a di(C_1 - C_6) alkylaminocarbonyl group whose (C_1-C_6) alkyl groups may be the same or different; a mono (C_1-C_6) alkylaminothiocarbonyl group; a $di(C_1-C_6)$ alkylaminothiocarbonyl group

whose (C_1-C_6) alkyl groups may be the same or different; a $di(C_1-C_6)$ alkoxyphosphoryl group whose (C_1-C_6) alkoxy groups may be the same or different; a $di(C_1-C_6)$ alkoxythiophosphoryl group whose (C_1-C_6) alkoxy groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1 - C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenyl (C_1-C_6) C₄)alkyl group; a substituted phenyl(C₁-C₄)alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl

groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and r is an integer of 1 to 4),

provided that $\ensuremath{R^1}$ and $\ensuremath{R^2}$ are not hydrogen atoms at the same time,

 ${\sf R}^1$ and ${\sf R}^2$ may form a 4 to 7 membered ring by combining to each other, in which the ring may contain the same or different 1 to 3 hetero atoms selected from the group consisting of oxygen atom, sulfur atom and nitrogen atom,

Het is a heterocyclic group represented by any of the following formulas Q1 to Q22:

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$$0 = 5$$

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$$Q9 = \begin{array}{c} Xn & (O) p \\ Xn & 1 & 3 \\ (O) q & 6 & 5 \end{array} \qquad Q10 = \begin{array}{c} Xn & (O) p \\ Xn & 1 \\ (N) & 2 \\ (N) & 3 \\ (O) q & & & \\ \end{array} \qquad Q11 = \begin{array}{c} Xn & 4 & 3 \\ (N) & 2 & 3 \\ (N) & 2 & 3 \\ (N) & 3 & 3 \\ (O) & q & & \\ \end{array}$$

Q21=
$$\frac{Xn}{2N}$$
 $\frac{3}{N}$ $\frac{4}{5}$ Q22= $\frac{Xn}{2N}$ $\frac{3}{N}$ $\frac{4}{5}$

(wherein X, which may be the same or different, are halogen atoms; cyano groups; nitro groups; (C_3-C_6) cycloalkyl groups; halo(C_3-C_6)cycloalkyl groups; tri($C_1 C_6$) alkylsilyl groups whose (C_1-C_6) alkyl groups may be the same or different; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1 - C_6) alkylthio groups, (C_1 - C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo(C_1 - C_6)alkylsulfonyl groups; or $-A^3-R^8$ [wherein A^3 is -O-, -S-, -SO-, $-SO_2-$, $-N(R^6)$ - (wherein R^6 is as defined above), -C(=0) -, -C(=NOR 7) - (wherein R 7 is as defined above), a (C₁-C₆)alkylene group, a halo(C_1-C_6)alkylene group, a ($C_2 C_6$) alkenylene group, a halo (C_2-C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R⁸ is as follows:

(1) when A^3 is -O-, -S-, -SO-, -SO₂- or -N(R^6)- (wherein

 R^6 is as defined above), then R^8 is a halo(C_3 - C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkenyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1$ - C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^4-R^9$ (wherein A^4 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) C_6) alkylene group, a (C_3-C_6) alkenylene group, a halo (C_3-C_6) C_6) alkenylene group, a (C_3-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R^9 is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups,

halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^5-R^{10}$ (wherein ${\rm A^5}$ is -O-, -S-, -SO-, -SO $_2-$ or -C(=0), and ${\rm R^{10}}$ is a (C_1-C_6) alkyl group; a halo (C_1-C_6) alkyl group; a (C_3-C_6) C_6) alkenyl group; a halo(C_3-C_6) alkenyl group; a (C_3-C_6) cycloalkyl group; a halo(C3-C6)cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups)), when A^3 is -C(=O) - or $-C(=NOR^7)$ - (wherein R^7 is as defined above), then R^8 is a hydrogen atom; a $(C_1 C_6$) alkyl group; a halo (C_1-C_6) alkyl group; a (C_2-C_6) alkenyl group; a halo(C_2-C_6) alkenyl group; a (C_3-C_6)-

cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) C_6) alkoxy group; a (C_1-C_6) alkylthio group; a mono (C_1-C_6) alkylamino group; a di (C_1-C_6) alkylamino group whose (C_1-C_6) C_6) alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) $C_6)$ alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenylamino group; a substituted phenylamino group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1 - C_6)alkylthio groups, $halo(C_1-C_6)$ alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps} \texttt{halo}\,(C_1 - C_6)\,\texttt{alkylsulfinyl} \ \texttt{groups,} \ (C_1 - C_6)\,\texttt{alkylsulfonyl}$ groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo(C_1 - C_6)alkylthio groups, (C_1 - C_6)alkylsulfinyl groups, halo(C_1-C_6)alkylsulfinyl groups, (C_1-C_6)alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups, and

(3) when A^3 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) C_6) alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6) C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, then R^8 is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C₁-C₆)alkoxycarbonyl group; a $tri(C_1-C_6)$ alkylsilyl group whose (C_1-C_6) alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^6-R^{11}$ (wherein A^6 is -O-, -S-, -SOor $-SO_2-$, and R^{11} is a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) C_6) cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen

atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^7-R^{12}$ (wherein A^7 is a (C_1-C_6) alkylene group, a halo(C_1-C_6)alkylene group, a (C_2-C_6)alkenylene group, a halo (C_2-C_6) alkenylene group, a (C_2-C_6) C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R^{12} is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) cycloalkyl group; a (C_1-C_6) alkoxy group; a halo(C_1-C_6) alkoxy group; a ($C_1 C_6$) alkylthio group; a halo (C_1-C_6) alkylthio group; a (C_1-C_6) C_6) alkylsulfinyl group; a halo (C_1-C_6) alkylsulfinyl group; a (C_1-C_6) alkylsulfonyl group; a halo (C_1-C_6) alkylsulfonyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio

groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenoxy group; a substituted phenoxy group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1 - C_6) alkyl groups, (C_1 - C_6) alkoxy groups, halo(C_1 - C_6) alkoxy groups, (C_1 - C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a phenylthio group; a substituted phenylthio group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, $\label{eq:coups} \verb|halo|(C_1-C_6)| alkylthio| groups, (C_1-C_6)| alkylsulfinyl groups,$ halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a heterocyclic group; or a substituted heterocyclic group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo(C_1-C_6)alkylsulfonyl groups))], and n is an integer of 0 to 3,

X may form a condensed ring by combining together with the adjacent atoms in the heterocyclic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C1-C6) alkyl groups; halo(C_1-C_6) alkyl groups; (C_1-C_6) alkoxy groups; halo(C_1 - C_6) alkoxy groups; (C_1-C_6) alkylthio groups; halo (C_1-C_6) C_6) alkylthio groups; (C_1-C_6) alkylsulfinyl groups; halo (C_1-C_6) alkylsulfinyl groups; (C_1-C_6) alkylsulfonyl groups; halo(C₁-C₆)alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C1-C6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo(C1-C6)alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $halo(C_1-C_6)$ alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups,

W is O, S or $N-R^{13}$ (wherein R^{13} is a $(C_1-C_6)-$ alkyl group; a halo (C_1-C_6) alkyl group; a (C_3-C_6) alkenyl

group; a halo(C_3-C_6)alkenyl group; a (C_3-C_6)alkynyl group; a halo(C_3-C_6) alkynyl group; a (C_1-C_6) alkoxy group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, $halo(C_1-C_6)$ alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps} \verb|halo|(C_1-C_6)| alkylsulfinyl groups, (C_1-C_6)| alkylsulfonyl \\$ groups and halo (C_1-C_6) alkylsulfonyl groups; a phenyl (C_1-C_6) C_6) alkyl group; or a substituted phenyl(C_1 - C_6) alkyl group having on the ring one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and p and q, which may be the same or different, are integers of 0 to 1),

 B^1 , B^2 , B^3 and B^4 , which may be the same or different, are carbon atoms or nitrogen atoms,

Y, which may be the same or different, are halogen atoms; cyano groups; nitro groups; halo (C_3-C_6) - cycloalkyl groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) -

alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups; sulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ (wherein A^3 and R^8 are as defined above), and m is an integer of 1 to 5,

Y may form a condensed ring by combining together with the adjacent carbon atoms in the aromatic ring, and said condensed ring may have one or more substituents, which may be the same or different, and are selected from halogen atoms; (C_1-C_6) alkyl groups; halo (C_1-C_6) alkyl groups; (C_1-C_6) alkoxy groups; halo (C_1-C_6) alkoxy groups; (C_1-C_6) alkylthio groups; halo (C_1-C_6) alkylthio groups; (C_1-C_6) alkylsulfinyl groups; halo (C_1-C_6) alkylsulfinyl groups; (C_1-C_6) alkylsulfonyl groups; halo (C_1-C_6) alkylsulfonyl groups; phenyl group; substituted phenyl groups having one or more substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups,

 C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; heterocyclic groups; and substituted heterocyclic groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, groups and halo (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups, and each of (C_1-C_6) alkylsulfonyl groups, and each of

provided that:

(1) when Het is Q2, Q6, Q7 or Q9 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 3-chloro-2-methyl group, 3-chloro-2,6-diethyl group, 5-chloro-2-methyl group, 2,6-diethyl group, 4-chloro-2-fluoro group and 2-ethyl-6-methyl group, (2) when Het is Q4 and B¹, B², B³ and B⁴ are carbon atoms at the same time, then Ym is other than 2,5-dichloro group, 2,4-difluoro group, 2,6-difluoro group, 3-chloro-2-methyl group, 5-chloro-2-methyl group, 5-fluoro-2-methyl group, 2,6-diethyl group, 2-ethyl-6-methyl group, 2-methoxy-5-nitro group, 2-methoxy-5-methyl group, 2,6-diethoxy group, 3-bromo-2-methyl group, 3-fluoro-2-methyl group, 3-iodo-2-methyl group, 3-cyano-2-methyl group, 3-difluoro-

methoxy-2-methyl group, 5-chloro-2-ethyl group, 2,5-dimethyl group, 2,3-dichloro group, 3-chloro-2,6-diethyl group, 4-trifluoromethyl group, 3-methoxy-carbonyl-2-methyl group, 3-trifluoromethyl-2-methyl group, 3,5-dichloro-2,6-diethyl group, 3,4-dichloro group, 3-(methoxycarbonylmethyloxy)-2-methyl group, 2-methyl-3-nitro group and 4-trifluoromethoxy group,

- (3) when Het is Q9, R^2 and R^3 are hydrogen atoms at the same time, Xn is a 2-phenyl group, R^1 is a n-propyl group or an i-propyl group and B^1 , B^2 , B^3 and B^4 are carbon atoms at the same time, then Ym is other than 4-pentafluoroethyl-2-methyl group, and
- (4) when Het is Q10 and B^1 , B^2 , B^3 and B^4 are carbon atoms at the same time, then Ym is other than 5-chloro-2-methyl group, 5-fluoro-2-methyl group and 2,5-dimethyl group).
- 2. A heterocyclic dicarboxylic acid diamide derivative according to claim 1, wherein Het is Q1, Q2, Q3 or Q4, R^1 is a (C_3-C_6) cycloalkyl group, a halo (C_3-C_6) cycloalkyl group or $-A^1-(R^4)$ r (wherein A^1 is a (C_1-C_8) -alkylene group, R^4 , which may be the same or different, are hydrogen atoms; halogen atoms; cyano groups; nitro groups; (C_1-C_6) alkoxycarbonyl groups; di (C_1-C_6) alkoxyphosphoryl groups whose (C_1-C_6) alkoxy groups may be the same or different; di (C_1-C_6) alkoxythiophosphoryl groups whose (C_1-C_6) alkoxy groups may be the same or different; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different

and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1 - C_6)alkoxy groups, (C_1 - C_6)alkylthio groups, $halo(C_1-C_6)$ alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps} \texttt{halo}\,(C_1 - C_6)\,\texttt{alkylsulfinyl} \ \texttt{groups,} \ (C_1 - C_6)\,\texttt{alkylsulfonyl}$ groups and halo (C_1-C_6) alkylsulfonyl groups; pyridyl groups; substituted pyridyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo(C_1-C_6) alkylsulfonyl groups; or $-A^2-R^5$ (wherein A^2 is -O-, -S-, -SO-, -SO₂-, -N(\mathbb{R}^6)- (wherein \mathbb{R}^6 is a hydrogen atom, a (C_1-C_6) alkylcarbonyl group, a halo (C_1-C_6) alkylcarbonyl group or a (C_1-C_6) alkoxycarbonyl group), or $-C(=NOR^7)$ - (wherein R^7 is a hydrogen atom, a $(C_1 C_6$) alkyl group or a halo (C_1-C_6) alkyl group), and R^5 is a hydrogen atom; a (C_1-C_6) alkyl group; a halo (C_1-C_6) alkyl group; a (C_1-C_6) cycloalkyl group; a halo (C_1-C_6) cycloalkyl group; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo(C_1-C_6)alkylsulfinyl groups, (C_1-C_6)alkyl-

sulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; a pyridyl group; or a substituted pyridyl group having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups), and r is an integer of 1 to 4), R^2 and R^3 , which may be the same or different, are hydrogen atoms or (C_1-C_6) alkyl groups, X and Y, which may be the same or different, are halogen atoms; cyano groups; nitro groups; phenyl groups; substituted phenyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, $(C_1 C_6$) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl groups and halo (C_1-C_6) alkylsulfonyl groups; pyridyl groups; substituted pyridyl groups having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, $halo(C_1-C_6)$ alkylthio groups, (C_1-C_6) alkylsulfinyl groups, halo (C_1-C_6) alkylsulfinyl groups, (C_1-C_6) alkylsulfonyl

groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^3-R^8$ [wherein A^3 is -O-, -S-, -SO-, $-SO_2-$, a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R^8 is as follows:

when A^3 is -O-, -S-, -SO- or -SO₂-, then R^8 is a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps_control} \mbox{halo}\,(\mbox{C_1-$$$$}\mbox{C_6})\,\mbox{alkylsulfinyl groups, } (\mbox{C_1-$$$$$}\mbox{C_6})\,\mbox{alkylsulfinyl}$ groups and halo (C_1-C_6) alkylsulfonyl groups; a pyridyl group; a substituted pyridyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1 - C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps} \verb|halo|(C_1-C_6)| alkyl sulfinyl groups, (C_1-C_6)| alkyl sulfonyl \\$ groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^4-R^9$ (wherein A^4 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) alkylene group, a (C_3-C_6) alkenylene group, a halo (C_3-C_6) C_6) alkenylene group, a (C_3-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and R^9 is a hydrogen atom, a halogen atom or $-A^5-R^{10}$ (wherein A^5 is -O-, -S-, -SO- or

 $-SO_2-$, and R^{10} is a (C_1-C_6) alkyl group, a halo (C_1-C_6) alkyl group, a (C_3-C_6) alkenyl group, a halo (C_3-C_6) alkenyl group, a (C_3-C_6) cycloalkyl group or a halo (C_3-C_6) cycloalkyl group), and

(2) when A^3 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) C_6) alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6) C_6) alkenylene group, a (C_2-C_6) alkynylene group or a halo(C_3-C_6)alkynylene group, then R^8 is a hydrogen atom; a halogen atom; a (C_3-C_6) cycloalkyl group; a halo (C_3-C_6) C_6) cycloalkyl group; a tri(C_1 - C_6) alkylsilyl group whose (C_1-C_6) alkyl groups may be the same or different; a phenyl group; a substituted phenyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, halo(C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo(C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, $\label{eq:coups} \verb|halo|(C_1-C_6)| alkylthio| groups, (C_1-C_6)| alkylsulfinyl groups,$ $\label{eq:comps} \texttt{halo}\,(\texttt{C}_1-\texttt{C}_6)\,\texttt{alkylsulfinyl}\,\,\texttt{groups,}\,\,\,(\texttt{C}_1-\texttt{C}_6)\,\texttt{alkylsulfonyl}$ groups and halo (C_1-C_6) alkylsulfonyl groups; a pyridyl group; a substituted pyridyl group having one or more substituents which may be the same or different and are selected from halogen atoms, (C_1-C_6) alkyl groups, $halo(C_1-C_6)$ alkyl groups, (C_1-C_6) alkoxy groups, $halo(C_1-C_6)$ C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) C_6) alkylthio groups, (C_1-C_6) alkylsulfinyl groups, $\label{eq:comps} \verb|halo(C_1-C_6)| alkyl sulfinyl groups, (C_1-C_6)| alkyl sulfonyl \\$ groups and halo (C_1-C_6) alkylsulfonyl groups; or $-A^6-R^{11}$ (wherein A^6 is -O-, -S-, -SO- or -SO₂-, and R^{11} is $-A^7-R^{12}$

(wherein A^7 is a (C_1-C_6) alkylene group, a halo (C_1-C_6) - alkylene group, a (C_2-C_6) alkenylene group, a halo (C_2-C_6) - alkenylene group, a (C_2-C_6) alkynylene group or a halo (C_3-C_6) alkynylene group, and (C_3-C_6) alkynylene group, and (C_1-C_6) alkoxy group, a halo (C_1-C_6) alkoxy group, a halo (C_1-C_6) alkylthio group, a halo (C_1-C_6) alkylthio group, a (C_1-C_6) alkylsulfinyl group, a halo (C_1-C_6) alkylsulfinyl group or a halo (C_1-C_6) alkylsulfonyl group or a halo (C_1-C_6) alkylsulfonyl group or a halo (C_1-C_6) alkylsulfonyl group)), both (C_1-C_6) alkylsulfonyl group), both (C_1-C_6) alkylsulfonyl group), a halo (C_1-C_6) alkylsulfonyl group), and (C_1-C_6) alkylsulfonyl group), and (C_1-C_6) alkylsulfonyl group, and (C_1-C_6) alkylsu

- 3. A heterocyclic dicarboxylic acid diamide derivative according to claim 2, wherein X, which may be the same or different, are halogen atoms, nitro groups, halo (C_1-C_6) alkyl groups, halo (C_1-C_6) alkylthio groups, and n is an integer of 0 to 3.
- 4. A heterocyclic dicarboxylic acid diamide derivative according to claim 3, wherein Y, which may be the same or different, are halogen atoms, (C_1-C_6) alkyl groups, halo (C_1-C_6) alkyl groups, (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxy groups, (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxyhalo (C_1-C_6) alkoxy groups, halo (C_1-C_6) alkoxyhalo or halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylthio groups, halo (C_1-C_6) alkylsulfinyl groups or halo (C_1-C_6) alkylsulfonyl groups, and m is an integer of 1 to 5.

- 5. A heterocyclic dicarboxylic acid diamide derivative according to claim 4, wherein R^1 is a (C_1-C_6) alkyl group, a (C_1-C_6) alkoxy (C_1-C_8) alkyl group, a (C_1-C_6) alkylthio (C_1-C_8) alkyl group, a (C_1-C_6) alkyl-sulfinyl (C_1-C_8) alkyl group or a (C_1-C_6) alkylsulfonyl (C_1-C_8) alkyl group, and R^2 and R^3 , which may be the same or different, are hydrogen atoms or methyl groups.
- An agricultural and horticultural insecticide characterized by containing a heterocyclic dicarboxylic acid diamide derivative according to any one of claims 1 to 5 as an active ingredient.
- 7. A method for applying an agricultural and horticultural insecticide, characterized by treating a crop to be protected or soil with an effective amount of an agricultural and horticultural insecticide according to claim 6 in order to protect useful crops against insect pests.

ABSTRACT

Heterocyclic dicarboxylic acid diamide derivative represented by the general formula (I):

$$\begin{array}{c|c} Xn & Z^{1} \\ \hline NR^{1}R^{2} \\ \hline R_{3} & B^{1} = B^{2} \\ \hline Z^{2} & R_{3} & B^{4} = B^{3} \end{array}$$
 (I)

wherein R^1 , R^2 and R^3 represent each H, optionally halogenated C_{3-6} cycloalkyl, etc.; Het represents a 5- or 6-membered heterocycle; X and Y represent each halocyano, nitro, optionally halogenated C_{3-6} cycloalkyl, optionally substituted phenyl, an optionally substituted phenyl, an optionally substituted heterocycle, etc; n is from 0 to 3; m is from 1 to 5; Z^1 and Z^2 represent each 0 or S; and B^1 to B^4 represent each C or N. Agricultural/horticultural insecticides having an excellent controlling effect on pest insects such as diamond-back moth (Pluntella xylostella) and tobacco cutworm (Spodoptera litura).

#8

RULE 63 (37 C.F.R. 1.63)

DECLARATION AND POWER OF ATTORNEY FOR UTILITY OR DESIGN PATENT APPLICATION IN THE UNITED STATES PATENT AND TRADEMARK OFFICE [] Declaration Submitted with Initial Filing or [] Declaration Submitted after Initial Filing (surcharge 37 CFR 1.16 (e) required)

	,
As a below named inventor, I hereby declare that my residence, post office address and citizenship are as stated below next to moriginal, first and sole inventor (if only one name is listed below) or an original, first and joint inventor (if plural names are listed which is claimed and for which a patent is sought on the INVENTION ENTITLED	ny name, and I believe I am the d below) of the subject matter
"HETEROCYCLIC DICARBOXYLIC ACID DIAMIDE DERIVATIVES, AGRICULTURAL AND	
[X] xttachcalcharate was filed on December 19, 2001, as U.S. Serial No. 10/018,46	specification of which is:
OR	
[X] was filed on (MM/DD/YYYY) <u>June 23, 2000</u> As United States Application Number (Attorney Docket No. PCT International Application No. <u>PCT/JP00/04136</u> and was amended on (if applicable)	
(II application No. FC170F00704136 and was amended on (II application)
I hereby state that I have reviewed and understand the contents of the above identified specification, including the claims, as a referred to above. I acknowledge the duty to disclose all information known to me to be material to patentability as defined in a continuation-in-part application, material information which becomes available between the filing date of the prior application international filing date of the continuation-in-part application.	37 C.F.R. 1.56 including for
I hereby claim foreign priority benefits under 35 U.S.C. 119(a) -(d) or 365(b) of any foreign application(s) for patent or inventor PCT international Application which designated at least one country other than the United States of America, listed below and checking the box, any foreign application for patent or inventors certificate, or any PCT international application having a filing on which priority is claimed.	have also identified below, by
PRIOR FOREIGN APPLICATION(S) Priority Not Claimed Certified Copy At	ttached?
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and further that these statements were made with the knowledge that willful false statements and the like so made are punishab both, under Section 1001 of Title 18 of the United States Code and that such willful false statements may jeopardize the validity issued thereon. I hereby appoint the registered practitioners represented by Customer No.: 20736 to prosecute this application and transact and tran	y of the application or any paten all business in the U.S. Patent
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